

2000 Annual Interim Sanitary Landfill Groundwater Monitoring Report

RECORDS ADMINISTRATION



APGB

by

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Savannah River Site

Aiken, South Carolina 29808

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INTERIM SANITARY LANDFILL GROUNDWATER MONITORING REPORT (U)

2000 ANNUAL

Publication Date: JANUARY 2001

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/ Reviewing Official

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Westinghouse Savannah River Company
Savannah River Site
Aiken, SC 29808

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**INTERIM SANITARY LANDFILL
2000 ANNUAL GROUNDWATER
MONITORING REPORT (U)
WSRC-TR-2000-000464**

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Environmental Protection Department

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INTRODUCTION

The SRS Interim Sanitary Landfill (ISL) opened in mid-1992 and operated until October 1999 under Domestic Waste Permit # 025500-1102. The ISL was subsequently closed and officially began the 30 year post-closure care period as of September 9, 2000, which is the effective date of the modified permit. This annual monitoring report was prepared in accordance with the modified permit and the SCDHEC approved *Closure/Post-Closure Plan, Q-CLP-G-00008, Rev. 2, October 26, 1998*.

Fifteen wells shown in Figure 1 were each sampled semi-annually during Calendar Year 2000 in accordance with *Closure/Post-Closure Plan, Q-CLP-G-00008, Rev. 2, October 26, 1998*. Sampling was performed during the first and third quarters of 2000. Sampling and analysis for the Appendix II list was performed in first quarter, and again in the third quarter for any analytes that were detected above background. Sampling and analysis for the Appendix I list was performed during third quarter 2000. The list of analytes is presented in Appendix A, Groundwater Protection Standards. The analytical results appear in Appendix B for the First and Second Halves of 2000. The vadose zone methane monitoring data is presented and discussed in Appendix D.

The well sampling and analyses were conducted in accordance with WSRC Procedure Manual 3Q5, Hydrogeologic Data Collection.

The report includes a discussion of the groundwater flow direction and rate, the groundwater analytical results, and the methane monitoring results. A key to reading the data tables in the preface of Appendix B. Figures are provided in Appendix C.

DISCUSSION

Flow Direction and Rate

The uppermost aquifer beneath the ISL is referred to as the Steed Pond Aquifer. All previous studies indicate that flow in the Steed Pond was to the southeast. Potentiometric data from the wells in the vicinity of the ISL are generally consistent with the regional flow. Figures 2, 3 and 4 show the potentiometric surface for three quarters spanning over a year. The flow direction is consistent over this period, showing a southerly and east-southeasterly component (relative to true north) that may be related to increased recharge over the ISL. The increased recharge would result from the previous condition of the waste cell as a low excavated area that would collect runoff. The closure of the ISL and associated cover placement over the waste cell may influence this observed flow pattern over time.

Flow rate can be estimated using the following equation:

$$\text{Flow (ft/day)} = \frac{\text{Hydraulic Conductivity (ft/day)} \times \text{dh(ft)}}{\text{Porosity (unitless)} \times \text{dl(ft)}}$$

where the hydraulic conductivity is assumed to be 16 ft/day, the effective porosity value is 20 percent, the change in head is dh, and the horizontal distance along each flow arrow is dl.

The flow rate was estimated in two directions indicated by flow arrows A-A' and B-B' on Figures 3 and 4.

Based on first quarter 2000 potentiometric data (Figure 3), the calculations are as follows:

Line A-A'

$$(16/.20) \times (4 \text{ ft}/261.4 \text{ ft}) = 1.22 \text{ ft/day or } 446.85 \text{ ft/year}$$

Line B-B'

$$16/.20 \times 4 \text{ ft}/603.5 \text{ ft} = 0.53 \text{ ft/day or } 193.5 \text{ ft/year}$$

Based on third quarter 2000 potentiometric data (Figure 4), the calculations are as follows:

Line A-A'

$$16/.20 \times 4 \text{ ft}/273.6 \text{ ft} = 1.16 \text{ ft/day or } 426.9 \text{ ft/year}$$

Line B-B'

$$16/.20 \times 4 \text{ ft}/668.7 \text{ ft} = 0.48 \text{ ft/day or } 174.7 \text{ ft/year}$$

Analytical Results

Groundwater Monitoring

The analytical data for the first and third quarter groundwater sampling events are provided in Appendix B. Results for analytes that exceeded the Groundwater Protection Standards (GWPS) shown in Table A.1 are discussed in the following text.

During first quarter 2000, five wells had analytes detected at levels exceeding GWPSs.

- Mercury (total recoverable) was elevated in well LFW-34 with a maximum concentration of 3.49 ug/L. However, mercury was not detected in the next sampling event (third quarter 2000).
- Trichlorofluoromethane was elevated in three wells; LFW-34 with a maximum concentration of 490 ug/L, LFW 32 with a concentration of 160 ug/L, and LFW 74D with a concentration of 21 ug/L.
- Zinc was elevated in two wells above the GWPS tolerance limit of 29.3 ug/L Appendix I constituents; LFW 76 with a maximum concentration of 35.6 ug/L, and LFW 78 with a maximum concentration of 32.7 ug/L.

During third quarter 2000, five wells had analytes detected at levels exceeding GWPSs.

- Copper (total recoverable) was elevated above the GWPS tolerance limit of 29.8 ug/L, Appendix I constituents in well LFW 31, with a maximum concentration of 40.8 ug/L. However, this concentration is less than the Primary Drinking Water Standard of 1000 ug/L.
- Dichloromethane (methylene chloride) was elevated in well LFW 43B, with a maximum concentration of 8.32 ug/L. This is a background well.
- Specific Conductance was elevated above the GWPS tolerance limits for Appendix I constituents in well LFW 74D, with a value of 48.5 uS/cm.
- Thallium (total recoverable) was elevated in well LFW 43D, with a maximum concentration of 10.1 ug/L. This is a background well.
- Trichlorofluoromethane was elevated in well LFW 34, with a maximum concentration of 450 ug/L.
- Zinc was elevated above the GWPS tolerance limit of 29.3 ug/L, Appendix I in well LFW 31, with a maximum concentration of 40.2 ug/L

Methane Monitoring

Results of the methane monitoring are provided in Appendix D, Table D.1. Methane was sampled at methane monitoring wells LGM 1, 2, 3 and 4. The results ranged from 0.51 vol% to 53.7 vol%. Methane levels exceeded the 5% LEL at LGM 1 and LGM 4.

In addition, extensive multilevel soil gas sampling was done around the Interim Sanitary Landfill to assess the extent to which methane in the soil had migrated outward from the landfill. The results were submitted to SCDHEC on September 28th, 2000 (letter, D.G. Wells to R.E. Schweitzer). The results of the soil gas sampling show that little or no migration has occurred. The submittal to SCDHEC also proposed fitting the existing methane monitoring wells with check valves.

CONCLUSION

Potentiometric data from the wells in the vicinity of the ISL are generally consistent with the southeast regional flow. Figures 2, 3, and 4 indicate a flow direction with southerly and east-southeasterly components (relative to true north), that may be related to increased recharge over the ISL. The closure of the ISL and associated cover placement over the waste cell may influence this observed flow pattern over time. Ground water flow rates range from 175 to 447 ft/year.

Mercury was detected above its GWPS of 2.0 ug/L (Final PDWS) in well LFW-34 downgradient of the ISL. Mercury is being monitored semiannually along with Appendix I analytes. Trichlorofluoromethane was detected above its GWPS of 20 ug/L (EPA Method 8240 Detection Limit) in three wells in the first quarter 2000 and one well (LFW 34) in the third quarter 2000. Well LFW 44D, which is about 500 feet downgradient from LFW 34, will be added to the assessment monitoring program. Assessment monitoring will continue for Zinc, Copper, Thallium, and Dichloromethane.

No new analytes were identified that were detected above the GWPS.

The vadose zone methane monitoring data and prior correspondence to the SCDHEC discussing the data is included in Appendix D.

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Appendix A

Groundwater Protection Standards

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Table A.1 Groundwater Protection Standards

| <u>CAS</u> | <u>Analyte Name</u> | <u>Synonym</u> | <u>Unit</u> | <u>Limit</u> | <u>Source†</u> |
|------------|--------------------------------|-----------------------------------|-------------|--------------|-------------------------|
| 630206 | 1,1,1,2-Tetrachloroethane | | µg/L | 20 | EPA Method 8240 |
| 71556 | 1,1,1-Trichloroethane | | µg/L | 200 | Final PDWS (EPA, 1998a) |
| 79345 | 1,1,2,2-Tetrachloroethane | | µg/L | 100 | EPA Method 8240 |
| 79005 | 1,1,2-Trichloroethane | | µg/L | 5 | Final PDWS (EPA, 1998a) |
| 75343 | 1,1-Dichloroethane | | µg/L | 20 | EPA Method 8240 |
| 75354 | 1,1-Dichloroethene | | µg/L | 7 | Final PDWS (EPA, 1998a) |
| 563586 | 1,1-Dichloropropene | | ug/L | 5 | EPA Method 8260B |
| 96184 | 1,2,3-Trichloropropane | | µg/L | 20 | EPA Method 8240 |
| 95943 | 1,2,4,5-Tetrachlorobenzene | | µg/L | 162 | EPA Method 8270 |
| 120821 | 1,2,4-Trichlorobenzene | | µg/L | 70 | Final PDWS (EPA, 1998a) |
| 96128 | 1,2-Dibromo-3-chloropropane | | µg/L | 0.2 | Final PDWS (EPA, 1998a) |
| 106934 | 1,2-Dibromoethane | | µg/L | 0.05 | Final PDWS (EPA, 1998a) |
| 95501 | 1,2-Dichlorobenzene | o-Dichlorobenzene | µg/L | 600 | Final PDWS (EPA, 1998a) |
| 107062 | 1,2-Dichloroethane | | µg/L | 5 | Final PDWS (EPA, 1998a) |
| 78875 | 1,2-Dichloropropane | | µg/L | 5 | Final PDWS (EPA, 1998a) |
| 99354 | 1,3,5-Trinitrobenzene | sym-Trinitrobenzene | µg/L | 162 | EPA Method 8270 |
| 541731 | 1,3-Dichlorobenzene | m-Dichlorobenzene | µg/L | 162 | EPA Method 8270 |
| 142289 | 1,3-Dichloropropane | | ug/L | 5 | EPAMethod 8260B |
| 99650 | 1,3-Dinitrobenzene | m-Dinitrobenzene | µg/L | 162 | EPA Method 8270 |
| 106467 | 1,4-Dichlorobenzene | | µg/L | 75 | Final PDWS (EPA, 1998a) |
| 130154 | 1,4-Naphthoquinone | p-Dichlorobenzene | µg/L | 162 | EPA Method 8270 |
| 134327 | 1-Naphthylamine | | µg/L | 162 | EPA Method 8270 |
| 594207 | 2,2-Dichloropropane | | µg/L | 10 | EPA Method 8260 |
| 108601 | 2,2-Oxybis(1-chloropropane) | Picloram | µg/L | 500 | Final PDWS (EPA, 1998a) |
| 58902 | 2,3,4,6-Tetrachlorophenol | | µg/L | 167 | EPA Method 8270 |
| 93765 | 2,4,5-T | 2,4,5-Trichlorophenoxyacetic acid | µg/L | 0.5 | EPA Method 8150 |
| 95954 | 2,4,5-Trichlorophenol | | µg/L | 10 | EPA Method 8270 |
| 88062 | 2,4,6-Trichlorophenol | | µg/L | 1 | EPA Method 8270 |
| 120832 | 2,4-Dichlorophenol | | µg/L | 10.2 | EPA Method 8270 |
| 94757 | 2,4-Dichlorophenoxyacetic acid | 2,4-D | µg/L | 70 | Final PDWS (EPA, 1998a) |
| 105679 | 2,4-Dimethyl phenol | | µg/L | 10.2 | EPA Method 8270 |
| 51285 | 2,4-Dinitrophenol | | µg/L | 102 | EPA Method 8270 |
| 121142 | 2,4-Dinitrotoluene | | µg/L | 1 | EPA Method 8270 |
| 87650 | 2,6-Dichlorophenol | | µg/L | 167 | EPA Method 8270 |
| 606202 | 2,6-Dinitrotoluene | | µg/L | 1 | EPA Method 8270 |
| 53963 | 2-Acetylaminofluorene | | µg/L | 162 | EPA Method 8270 |
| 91587 | 2-Chloronaphthalene | | µg/L | 10.2 | EPA Method 8240 |
| 95578 | 2-Chlorophenol | | µg/L | 10.2 | EPA Method 8270 |
| 591786 | 2-Hexanone | | µg/L | 100 | EPA Method 8240 |
| 534521 | 2-Methyl-4,6-dinitrophenol | 4,6-Dintro-o-cresol | µg/L | 102 | EPA Method 8270 |
| 91576 | 2-Methylnaphthalene | | µg/L | 10 | EPA Method 8270 |
| 91598 | 2-Naphthylamine | | µg/L | 162 | EPA Method 8270 |
| 88744 | 2-Nitroaniline | o-Nitroaniline | µg/L | 10 | EPA Method 8270 |
| 88755 | 2-Nitrophenol | o-Nitrophenol | ug/L | 10.2 | EPA Method 8270 |
| 88857 | 2-sec-Butyl-4,6-dinitrophenol | Dinoseb | ug/L | 7 | Final PDWS (EPA, 1998a) |
| 91941 | 3,3'-Dichlorobenzidine | | µg/L | 10.2 | EPA Method 8270 |
| 119937 | 3,3'-Dimethoxybenzidine | 3,3'-Dimethylbenzidine | µg/L | 162 | EPA Method 8270 |
| 56495 | 3-Methylcholanthrene | | µg/L | 162 | EPA Method 8270 |

| <u>CAS</u> | <u>Analyte Name</u> | <u>Synonym</u> | <u>Unit</u> | <u>Limit</u> | <u>Source†</u> |
|----------------|--------------------------------------|-----------------------------------|-------------|--------------|--------------------------------|
| 92671 | 4-Aminobiphenyl | a- Aminobiphenyl | µg/L | 162 | EPA Method 8270 |
| 101553 | 4-Bromophenyl phenyl ether | | µg/L | 10.2 | EPA Method 8270 |
| 106478 | 4-Chloroaniline | p-Chloroaniline | µg/L | 10 | EPA Method 8270 |
| 59507 | 4-Chloro-m-cresol | p-Cholo-m-cresol | µg/L | 10.2 | EPA Method 8270 |
| 7005723 | 4-Chlorophenyl phenyl ether | | µg/L | 10.2 | EPA Method 8270 |
| 100016 | 4-Nitroaniline | p-Nitroaniline | µg/L | 10 | EPA Method 8270 |
| 100027 | 4-Nitrophenol | p-Nitrophenol | µg/L | 10.2 | EPA Method 8270 |
| 99558 | 5-Nitro-o-toluidine | | µg/L | 162 | EPA Method 8270 |
| 57976 | 7,12-Dimethylbenz[a]anthracene | | µg/L | 162 | EPA Method 8270 |
| 83329 | Acenaphthene | | µg/L | 10.2 | EPA Method 8270 |
| 208968 | Acenaphthylene | | µg/L | 10.2 | EPA Method 8270 |
| 67641 | Acetone | | µg/L | 20 | EPA Method 8260B |
| 75058 | Acetonitrile (Methyl cyanide) | | µg/L | 100 | EPA Method 8240 |
| 98862 | Acetophenone | | µg/L | 170 | EPA Method 8270 |
| 107028 | Acrolein | | µg/L | 333 | EPA Method 8240 |
| 107131 | Acrylonitrile | | µg/L | 500 | EPA Method 8240 |
| 309002 | Aldrin | | µg/L | 0.8 | EPA Method 8080 |
| 107051 | Allyl chloride | | µg/L | 833 | EPA Method 8240 |
| 319846 | alpha-Benzene hexachloride | alpha-BHC | µg/L | 0.3 | EPA Method 8080 |
| 120127 | Anthracene | | µg/L | 10.2 | EPA Method 8270 |
| 7440360 | Antimony, total recoverable | | µg/L | 6 | Final PDWS (EPA, 1998a) |
| 7440382 | Arsenic, total recoverable | | µg/L | 50 | Final PDWS (EPA, 1998a) |
| 7440393 | Barium, total recoverable | | µg/L | 2000 | Final PDWS (EPA, 1998a) |
| 71432 | Benzene | | µg/L | 5 | Final PDWS (EPA, 1998a) |
| 56553 | Benzo[a]anthracene | | µg/L | 0.1 | Proposed PDWS (EPA, 1990) |
| 50328 | Benzo[a]pyrene | | µg/L | 0.2 | Final PDWS (EPA, 1998a) |
| 205992 | Benzo[b]fluoranthene | | µg/L | 0.2 | Proposed PDWS (EPA, 1990) |
| 191242 | Benzo[g,h,i]perylene | | µg/L | 10.2 | EPA Method 8270 |
| 207089 | Benzo[k]fluoranthene | | µg/L | 0.2 | Proposed PDWS (EPA, 1990) |
| 100516 | Benzyl alcohol | | µg/L | 10 | EPA Method 8270 |
| 7440417 | Beryllium, total recoverable | | µg/L | 4 | Final PDWS (EPA, 1998a) |
| 319857 | beta-Benzene hexachloride | beta-BHC | µg/L | 0.5 | EPA Method 8080 |
| 111911 | Bis(2-chloroethoxy) methane | bis[2-chloro-l-methylethyl] ether | µg/L | 10.2 | EPA Method 8270 |
| 111444 | Bis(2-chloroethyl) ether | | µg/L | 10.2 | EPA Method 8270 |
| 117817 | Bis(2-ethylhexyl) phthalate | | µg/L | 6 | Final PDWS (EPA, 1998a) |
| 74975 | Bromochloromethane | | µg/L | 10 | EPA Method 8260 |
| 75274 | Bromodichloromethane | | µg/L | 100 | Final PDWS (EPA, 1998a) |
| 75252 | Bromoform | | µg/L | 100 | Final PDWS (EPA, 1998a) |
| 74839 | Bromomethane | Methyl bromide | µg/L | 20 | EPA Method 8240 |
| 85687 | Butylbenzyl phthalate | | ug/L | 10 | Set by EPD/EMS |
| 7440439 | Cadmium, total recoverable | | µg/L | 5 | Final PDWS (EPA, 1998a) |
| 75150 | Carbon disulfide | | µg/L | 50 | EPA Method 8240 |
| 56235 | Carbon tetrachloride | | µg/L | 5 | Final PDWS (EPA, 1998a) |
| 57749 | Chlordane | | µg/L | 2 | Final PDWS (EPA, 1998a) |
| 108907 | Chlorobenzene | | µg/L | 100 | Final PDWS (EPA, 1998a) |
| 510156 | Chlorobenzilate | | µg/L | 162 | EPA Method 8270 |
| 75003 | Chloroethane | | µg/L | 20 | EPA Method 8240 |
| 75014 | Chloroethene (Vinyl chloride) | Vinyl chloride | µg/L | 2 | Final PDWS (EPA, 1998a) |
| 67663 | Chloroform | | µg/L | 100 | Final PDWS (EPA, 1998a) |
| 74873 | Chloromethane | (Methyl chloride) | µg/L | 20 | EPA Method 8240 |

| <u>CAS</u> | <u>Analyte Name</u> | <u>Synonym</u> | <u>Unit</u> | <u>Limit</u> | <u>Source†</u> |
|------------|-----------------------------|--------------------------|-------------|--------------|---------------------------|
| 126998 | Chloroprene | | µg/L | 3330 | EPA Method 8240 |
| 7440473 | Chromium, total recoverable | | µg/L | 100 | Final PDWS (EPA, 1998a) |
| 218019 | Chrysene | | µg/L | 0.2 | Proposed PDWS (EPA, 1990) |
| 156592 | cis-1,2-Dichloroethene | cis-1,2-Dichloroethylene | µg/L | 70 | Final PDWS (EPA, 1998a) |
| 10061015 | cis-1,3-Dichloropropene | | µg/L | 20 | EPA Method 8240 |
| 7440484 | Cobalt, total recoverable | | µg/L | 100 | EPA Method 6010 |
| 7440508 | Copper, total recoverable | | µg/L | 29.8 | EPA Method 6010 |
| 57125 | Cyanide | | µg/L | 200 | Final PDWS (EPA, 1998a) |
| 319868 | delta-Benzene hexachloride | delta-BHC | µg/L | 0.5 | EPA Method 8080 |
| 2303164 | Diallate | | µg/L | 162 | EPA Method 8270 |
| 53703 | Dibenz[a,h]anthracene | | µg/L | 0.3 | Proposed PDWS (EPA, 1990) |
| 132649 | Dibenzofuran | | µg/L | 10 | EPA Method 8270 |
| 124481 | Dibromochloromethane | | µg/L | 100 | Final PDWS (EPA, 1998a) |
| 74953 | Dibromomethane | Methylene bromide | ug/L | 20 | EPA Method 8240 |
| 75718 | Dichlorodifluoromethane | | µg/L | 20 | EPA Method 8240 |
| 75092 | Dichloromethane | Methylene Chloride | µg/L | 5 | Final PDWS (EPA, 1998a) |
| 60571 | Dieldrin | | µg/L | 8.3 | EPA Method 8080 |
| 84662 | Diethyl phthalate | | ug/L | 10 | EPA Method 8270C |
| 60515 | Dimethoate | | µg/L | 162 | EPA Method 8270 |
| 131113 | Dimethyl phthalate | | ug/L | 10 | EPA Method 8270C |
| 84742 | Di-n-butyl phthalate | | ug/L | 10 | EPA Method 8270C |
| 117840 | Di-n-octyl phthalate | | ug/L | 10 | EPA Method 8270C |
| 122394 | Diphenylamine | | µg/L | 162 | EPA Method 8270 |
| 298044 | Disulfoton | | µg/L | 162 | EPA Method 8270 |
| 959988 | Endosulfan I | | µg/L | 0.5 | EPA Method 8080 |
| 33213659 | Endosulfan II | | µg/L | 1.1 | EPA Method 8080 |
| 1031078 | Endosulfan sulfate | | µg/L | 1.1 | EPA Method 8080 |
| 72208 | Endrin | | µg/L | 2 | Final PDWS (EPA, 1998a) |
| 7421934 | Endrin aldehyde | | µg/L | 1.7 | EPA Method 8080 |
| 97632 | Ethyl methacrylate | | µg/L | 5 | EPA Method 8270 |
| 62500 | Ethyl methanesulfonate | | µg/L | 162 | EPA Method 8270 |
| 100414 | Ethylbenzene | | µg/L | 700 | Final PDWS (EPA, 1998a) |
| 52857 | Famphur | | µg/L | 162 | EPA Method 8270 |
| 206440 | Fluoranthene | | µg/L | 10.2 | EPA Method 8270 |
| 86737 | Fluorene | | µg/L | 10.2 | EPA Method 8270 |
| 76448 | Heptachlor | | µg/L | 0.4 | Final PDWS (EPA, 1998a) |
| 1024573 | Heptachlor epoxide | | µg/L | 0.2 | Final PDWS (EPA, 1998a) |
| 118741 | Hexachlorobenzene | | µg/L | 1 | Final PDWS (EPA, 1998a) |
| 87683 | Hexachlorobutadiene | | µg/L | 10 | EPA Method 8270 |
| 77474 | Hexachlorocyclopentadiene | | µg/L | 50 | Final PDWS (EPA, 1998a) |
| 67721 | Hexachloroethane | | µg/L | 1 | EPA Method 8270 |
| 1888717 | Hexachloropropene | | µg/L | 162 | EPA Method 8270 |
| 193395 | Indeno[1,2,3-c,d]pyrene | | µg/L | 1 | EPA Method 8270 |
| 74884 | Iodomethane (Methyl iodide) | | µg/L | 250 | EPA Method 8240 |
| 78831 | Isobutyl alcohol | | µg/L | 1669 | EPA Method 8240 |
| 465736 | Isodrin | | µg/L | 162 | EPA Method 8270 |
| 78591 | Isophorone | | µg/L | 10.2 | EPA Method 8270 |
| 120581 | Isosafrole | | µg/L | 162 | EPA Method 8270 |
| 143500 | Kepone | | µg/L | 162 | EPA Method 8270 |

| <u>CAS</u> | <u>Analyte Name</u> | <u>Synonym</u> | <u>Unit</u> | <u>Limit</u> | <u>Source†</u> |
|--|----------------------------------|-----------------------------|-------------|--------------|--------------------------------|
| Appendix I, per WSRC-RP-98-01263 | | | | | |
| 7439921 | Lead, total recoverable | | µg/L | 18 | |
| 58899 | Lindane | gamma-BHC | µg/L | 0.2 | Final PDWS (EPA, 1998a) |
| 108394 | m-Cresol (3-Methylphenol) | | µg/L | 100 | EPA Method 8270 |
| 7439976 | Mercury, total recoverable | | µg/L | 2 | Final PDWS (EPA, 1998a) |
| 126987 | Methacrylonitrile | | µg/L | 833 | EPA Method 8240 |
| 91805 | Methapyrilene | | µg/L | 162 | EPA Method 8270 |
| 72435 | Methoxychlor | | µg/L | 40 | Final PDWS (EPA, 1998a) |
| 78933 | Methyl ethyl ketone | 2-Butanone | ug/L | 20 | EPA Method 8260B |
| 108101 | Methyl isobutyl ketone | 4-methyl-2-pentanone | ug/L | 10 | EPA Method 8260B |
| 80626 | Methyl methacrylate | | µg/L | 100 | EPA Method 8270 |
| 66273 | Methyl methanesulfonate | | µg/L | 162 | EPA Method 8270 |
| 99092 | m-Nitroaniline | | µg/L | 10 | EPA Method 8270 |
| 108383 | m-Xylene | | µg/L | NA | NA |
| 91203 | Naphthalene | | µg/L | 167 | EPA Method 8270 |
| 7440020 | Nickel, total recoverable | | µg/L | 100 | Final PDWS (EPA, 1998a) |
| 98953 | Nitrobenzene | | µg/L | 10.2 | EPA Method 8270 |
| 55185 | N-Nitrosodiethylamine | | µg/L | 162 | EPA Method 8270 |
| 62759 | N-Nitrosodimethylamine | | µg/L | 167 | EPA Method 8270 |
| 924163 | N-Nitrosodi-n-butylamine | | µg/L | 162 | EPA Method 8270 |
| 86306 | N-Nitrosodiphenylamine | | µg/L | 10.2 | EPA Method 8270 |
| 621647 | N-Nitrosodipropylamine | | µg/L | 10.2 | EPA Method 8270 |
| 10595956 | N-Nitrosomethylethylamine | | µg/L | 162 | EPA Method 8270 |
| 100754 | N-Nitrosopiperidine | | µg/L | 162 | EPA Method 8270 |
| 930552 | N-Nitrosopyrrolidine | | µg/L | 162 | EPA Method 8270 |
| 126681 | O,O,O-Triethyl phosphorothioate | | µg/L | 162 | EPA Method 8270 |
| 95487 | o-Cresol (2-Methylphenol) | | µg/L | 10 | EPA Method 8270 |
| 95534 | o-Toluidine | | µg/L | 162 | EPA Method 8270 |
| 95476 | o-Xylene | | µg/L | 10 | EPA Method 8260 |
| 72548 | p,p'-DDD | 4,4-DDD | µg/L | 1.1 | EPA Method 8080 |
| 72559 | p,p'-DDE | 4,4-DDE | µg/L | 0.5 | EPA Method 8080 |
| 50293 | p,p'-DDT | 4,4-DDT | g/L | 1.7 | EPA Method 8080 |
| 56382 | Parathion | | µg/L | 0.8 | EPA Method 8080 |
| 298000 | Parathion methyl | | µg/L | 0.8 | EPA Method 8080 |
| 1336363 | PCBs | Polychlorinated Biphenyls | | NA | NA |
| 106445 | p-Cresol (4-Methylphenol) | | µg/L | 120 | EPA Method 8270 |
| 60117 | p-Dimethylaminoazobenzene | | µg/L | 162 | EPA Method 8270 |
| 608935 | Pentachlorobenzene | | µg/L | 162 | EPA Method 8270 |
| 82688 | Pentachloronitrobenzene | | µg/L | 162 | EPA Method 8270 |
| 87865 | Pentachlorophenol | | µg/L | 1 | Final PDWS (EPA, 1998a) |
| 4.11 - Appendix I, per WSRC-RP-98-01263 | | | | | |
| | pH | | pH | 6-15 | |
| 62442 | Phenacetin | | g/L | 162 | EPA Method 8270 |
| 85018 | Phenanthrene | | µg/L | 10.2 | EPA Method 8270 |
| 108952 | Phenol | | µg/L | 167 | EPA Method 8270 |
| 298022 | Phorate | | µg/L | 1.7 | EPA Method 8080 |
| 106503 | p-Phenylenediamine | | µg/L | 162 | EPA Method 8270 |
| 23950585 | Pronamid | | µg/L | 162 | EPA Method 8270 |
| 107120 | Propionitrile | | µg/L | 3330 | EPA Method 8240 |
| 106423 | p-Xylene | | | NA | NA |
| 129000 | Pyrene | | µg/L | 10.2 | EPA Method 8270 |

| <u>CAS</u> | <u>Analyte Name</u> | <u>Synonym</u> | <u>Unit</u> | <u>Limit</u> | <u>Source</u> [†] |
|-----------------|------------------------------------|--|-------------|--------------|--|
| 94597 | Safrole | | µg/L | 162 | EPA Method 8270 |
| 7782492 | Selenium, total recoverable | | µg/L | 50 | Final PDWS (EPA, 1998a) |
| 7440224 | Silver, total recoverable | | µg/L | 20 | EPA Method 6010B |
| 93721 | Silvex | 2,4,5-TP (Silvex) | µg/L | 50 | Final PDWS (EPA, 1998a) |
| | Specific Conductance | | µS/cm | 37.0 | 10.6 - Appendix I, per WSRC-RP-98-01263 |
| 100425 | Styrene | | µg/L | 100 | Final PDWS (EPA, 1998a) |
| 18496258 | Sulfide | | µg/L | 16700 | EPA Method 9030 |
| 127184 | Tetrachloroethylene | | µg/L | 5 | Final PDWS (EPA, 1998a) |
| 7440280 | Thallium, total recoverable | | µg/L | 2 | Final PDWS (EPA, 1998a) |
| | | 0,0-Diethyl 0-2-Pyrazinyl Phosphorothioate | | | |
| 297972 | Thionazin | | µg/L | 162 | EPA Method 8270 |
| 7440315 | Tin | | µg/L | 500 | EPA Method 282.2 |
| 108883 | Toluene | | µg/L | 1000 | Final PDWS (EPA, 1998a) |
| 8001352 | Toxaphene | | µg/L | 3 | Final PDWS (EPA, 1998a) |
| 156605 | trans-1,2-Dichloroethylene | | µg/L | 100 | Final PDWS (EPA, 1998a) |
| 10061026 | trans-1,3-Dichloropropene | | µg/L | 20 | EPA Method 8240 |
| 110576 | trans-1,4-Dichloro-2-butene | | µg/L | 500 | EPA Method 8240 |
| 79016 | Trichloroethylene | | µg/L | 5 | Final PDWS (EPA, 1998a) |
| 75694 | Trichlorofluoromethane | | µg/L | 20 | EPA Method 8240 |
| 7440622 | Vanadium | | µg/L | 133 | EPA Method 6010 |
| 108054 | Vinyl acetate | | µg/L | 100 | EPA Method 8240 |
| 1330207 | Xylenes | Xylene (Total) | µg/L | 10000 | Final PDWS (EPA 1998a) |
| | | | | | Appendix I, per WSRC-RP-98-01263 |
| 7440666 | Zinc, total recoverable | | µg/L | 29.3 | 98-01263 |

† Practical Quantitation Limits (PQLs) are provided as limits where EPA methods are listed as the "Source". References for dated sources are in the **References** section.

Note: Analytes from the Appendix I constituent list of the approved *Closure/Post-Closure Plan, Q-CLP-G-00008, Rev. 2, October 26, 1998* (WSRC-RP-98-01263) are shown in **bold text**. Analytes from the Appendix II list of the closure plan include all the analytes listed above with the exception of pH and specific conductance, which are only listed in Appendix I.

References

EPA (U.S. Environmental Protection Agency), 1977. **National Interim Primary Drinking Water Regulations**, EPA-570/9-76-003. Washington, DC.

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EPA (U.S. Environmental Protection Agency), 1991. *National Primary Drinking Water Regulations; Radionuclides; Proposed Rule*. **Federal Register**, July 18, 1991, pp. 33052-33127. Washington, DC.

EPA (U.S. Environmental Protection Agency), 1993a. *National Primary Drinking Water Regulations. Code of Federal Regulations*, Section 40, Part 141, pp. 592-732. Washington, DC.

EPA (U.S. Environmental Protection Agency), 1993b. *National Secondary Drinking Water Regulations. Code of Federal Regulations*, Section 40, Part 143, pp. 774-777. Washington, DC.

SCDHEC (South Carolina Department of Health and Environmental Control), 1981. **State Primary Drinking Water Regulations**, R.61-58.5. Columbia, SC.

EPA (U.S. Environmental Protection Agency), 1977. **National Interim Primary Drinking Water Regulations**, EPA-570/9-76-003. Washington, DC.

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SCDHEC (South Carolina Department of Health and Environmental Control), 1981. **State Primary Drinking Water Regulations**, R.61-58.5. Columbia, SC.

SRS (Westinghouse Savannah River Site), 1998. *Closure/Post-Closure Plan, Q-CLP-G-00008, Rev. 2, October 26, 1998 (WSRC-RP-98-01263)*. Aiken, SC.

Appendix B

Groundwater Monitoring Results Tables

Key to Reading the Tables

The following abbreviations may appear in the data tables:

Constituents

| | |
|---------------------|---|
| 1,2,3,4,6,7,8-HPCDD | 1,2,3,4,6,7,8-heptachlorodibenzo-p-dioxin |
| 1,2,3,4,6,7,8-HPCDF | 1,2,3,4,6,7,8-heptachlorodibenzo-p-furan |
| 1,2,3,4,7,8-HXCDD | 1,2,3,4,7,8-hexachlorodibenzo-p-dioxin |
| 1,2,3,4,7,8-HXCDF | 1,2,3,4,7,8-hexachlorodibenzo-p-furan |
| Lindane | gamma-benzene hexachloride |
| PCB | polychlorinated biphenyl |
| 1,2,3,7,8-PCDD | 1,2,3,7,8-pentachlorodibenzo-p-dioxin |
| 1,2,3,7,8-PCDF | 1,2,3,7,8-pentachlorodibenzo-p-furan |
| Sp. conductance | specific conductance |
| TCDD | tetrachlorodibenzo-p-dioxin |
| TCDF | tetrachlorodibenzo-p-furan |

Laboratories

| | |
|-----------|--|
| CN | Clemson Technical Center, Inc. |
| EM | Environmental Protection Department/Environmental Monitoring Section (EPD/EMS) Laboratory |
| GE and GP | General Engineering Laboratories, Inc. |
| SC | Savannah River Technology Center |
| SP | Spencer Testing Services, Inc. |
| TM | Thermo NUtech |
| WA and WS | Recra LabNet Philadelphia. |

Nomenclature

| | | |
|-----------|------|--|
| AZ | - | Aquifer Zone |
| CBAU | CBA | Crouch Branch Aquifer Unit (previously Black Creek) |
| CBCU | CBC | Crouch Branch Confining Unit |
| GAU | GAU | Gordon Aquifer Unit |
| GCCZ | GC | Green Clay Confining Zone |
| LAZ_UTRA | LAZ | Lower Aquifer Zone-Upper Three Runs Aquifer Unit |
| LLAZ | LLAZ | Lost Lake Aquifer Zone |
| LLLAZ | LL | Lower Lost Lake Aquifer Zone (previously Lower Congaree) |
| MAAZ | M | M-Area Aquifer Zone (previously Water Table) |
| MSAZ_CBCU | MCBC | Middle Sand Aquifer Zone-CBCU (previously Ellenton Sand) |
| PZ | PZ | Perched Zone |
| UAZ_UTRA | UAZ | Upper Aquifer Zone-Upper Three Runs Aquifer Unit |
| ULLAZ | UL | Upper Lost Lake Aquifer Zone (previously Upper Congaree) |
| UNKNOWN | UNK | Unknown |

Sampling Codes

| | |
|---|--|
| B | blank sample was collected |
| C | well was pumping continuously |
| D | well was dry |
| E | equipment blank was collected |
| I | well went dry during sampling; insufficient water to collect all samples |
| L | well went dry before sampling began; only depth to water can be determined |
| N | well was not stabilized before sampling began |
| P | inaccessibility or mechanical failure prevented sample collection and field analysis of the water |
| S | no water in standpipe |
| T | samples were collected, but some samples were not sent to the laboratory due to high turbidity |
| W | unable to sample well because of stabilization or sampling equipment failure; water-level measurements were obtained |
| X | well went dry during purging; samples collected after well recovered |

Sampling Methods

| | |
|---|---|
| B | sample collected using an open-bucket bailer |
| O | sample collected by method other than bailer or pump |
| P | sample collected using a bladder pump |
| S | sample collected using a single-speed centrifugal downhole pump |
| V | sample collected using a variable-speed pump |

Units

| | |
|--------|------------------------------|
| mg/L | milligrams per liter |
| msl | mean sea level |
| MSL | million structures per liter |
| NTU | nephelometric turbidity unit |
| pCi/L | picocuries per liter |
| pCi/mL | picocuries per milliliter |
| pH | pH unit |
| µg/L | micrograms per liter |
| µS/cm | microsiemens per centimeter |

Other Codes

| | |
|--------|--|
| CLP | EPA Functional Guideline Codes |
| CS | carbon steel |
| DF | <i>dilution factor</i> column in data tables |
| E | exponential notation (e.g., $1.1\text{E}-09 = 1.1 \times 10^{-9} = 0.0000000011$) |
| STORET | EPA STORET result qualifiers |
| Filt. | <i>Data results after application of the Data Usability filter</i> |
| GWPS | groundwater protection standard |
| MCS | monitoring constituents standard |
| Mod | <i>modifier</i> column in data tables |
| NDD | <i>"not decision data"</i> |
| PDWS | primary drinking water standard |
| PVC | polyvinyl chloride |
| ST | <i>exceeded the GWPS</i> column in data tables |
| TOC | top of casing |
| <EQL | less than the sample-specific estimated quantitation limit |

Results Below Detection

For radiological analyses, the analytical result field contains the result recorded on the analytical instrument and reported by the laboratory, even if it is negative. For nonradiological analyses, if the analyte is not detected, the sample-specific estimated quantitation limit (EQL) is entered into the result field and is reported with a less than [$<$] sign. The EQL is defined as the lowest concentration that can be achieved reliably within specified limits of precision and accuracy during routine laboratory operating conditions. The sample-specific EQL is modified for sample concentration or dilution or unusual aliquot size that affects analytical sensitivity.

Uncertainty and Data Usability

In April 1998, SCDHEC accepted guidance proposed by SRS to apply a method for minimizing uncertainty in compliance decisions potentially affecting long-term monitoring or remediation (letter, Taylor to Cook; April 21, 1998). The method is applied by processing or "filtering" the data, using the EPA Functional Guideline Codes applied by the laboratories to qualify the analytical results. By removing all data with a result qualifier of "L", "R", "U", and "J" from consideration, groundwater data users can ensure that only quantified numerical results are applied to the decision process. The output of the filtering process populates the "Filt" column as follows:

1) "Null" or "blank" – Data not remarked. The analytical result is acceptable for use as reported, and the result is not greater than an associated concentration limit for the analyte. If a concentration limit exists for the analyte, and the result is greater than the concentration limit, then the "Filt." Column will contain the applicable limit, and the "greater than" symbol. For example, in the case of trichloroethylene, ">5" would indicate the result exceeded a concentration limit of 5 ug/L.

Rationale: The best result would be one without qualifiers, so the preferred choice would be the maximum result that did not have any qualifiers.

2) "J", "L", "N", "NJ", or "JL" – "J" identifies that the analyte was positively identified; the associated numerical value is an estimated concentration of the analyte in the sample. "L" Indicates the sample result is off scale high. "JL" Indicates an estimated quantity of a sample that is off scale high. "N" is used for all TIC (tentatively identified compounds) and indicates the presence of an analyte for which there is presumptive evidence to make a tentative identification. "NJ" means the presence of an analyte that has been tentatively identified and the associated numerical value represents its approximate concentration.

Rationale: an estimate can still provide useful information. Although there may be a range of uncertainty around the actual value, the value itself may still grossly exceed a regulatory standard. However, a estimated value is less certain than an unqualified result. Therefore, this would be labeled as "NDD" (not decision data).

3) "U" - material analyzed for, but not detected. The analyte concentration is less than the sample specific Estimated Quantitation Limit and labeled "<EQL".

Rationale: a result above the detection limit would be chosen before a result below detection so that the process is not biased toward false negatives.

4) "UJ" - result is not above the reported sample quantitation limit, but the reported quantitation limit itself is approximate, and may not represent the actual limit of quantitation necessary to accurately and precisely measure the analyte in the sample.

Rationale: the additional qualifiers make this result less reliable for use than the "U" without qualifiers. These results would be labeled "<EQL".

5) "Rejected" – The sample results are rejected due to serious deficiencies in the ability to analyze the sample and meet quality control criteria. The presence or absence of the analyte cannot be verified.

Rationale: the only value in providing this result in the report is to indicate that the lab attempted to analyze the sample. If there are any other results available, the result with the "R" qualifier should not be reported. If it is reported, it is definitely "NDD" (not decision data).

Holding Times

Standard analytical methods include a limit, called holding time, on the maximum elapsed time between sample collection and extraction or analysis by the laboratory. In the data tables, the result qualifier Q in the "EPA" column indicates that holding time was exceeded. Analyses performed beyond holding times may not yield valid results.

The South Carolina Department of Health and Environmental Control (SCDHEC) allows only 15 minutes to elapse between sampling and analysis for pH. Thus, only field pH measurements can meet the holding time criterion; laboratory pH analyses always will exceed it.

The laboratory procedure used for the determination of specific conductance allows one day to elapse between sampling and analysis. Thus, laboratory specific conductance measurements may exceed the holding time criterion.

Data Qualification

The contract laboratories submit sample- or batch-specific quality assurance/quality control information either at the same time as analytical results or in a quarterly summary. Properly defined and used, data qualifiers can be a key component in assessing data usability. The EPA Functional Guideline Codes used by the analytical laboratories are shown in the CLP result qualifier column are defined below. These modifiers appear in the data tables under the column *CLP*. EPA STORET codes appear in the data tables under the column labeled *EPA*.

"CLP" Qualifiers - EPA Functional Guideline Codes

| | |
|---------|--|
| (Blank) | Data not remarked. The analytical result is acceptable for use as reported. |
| J | The analyte was positively identified; the associated numerical value is an estimated concentration of the analyte in the sample. |
| N | The analysis indicates the presence of an analyte for which there is presumptive evidence to make a tentative identification. Used for all TIC results. |
| R | The sample results are rejected due to serious deficiencies in the ability to analyze the sample and meet quality control criteria. The presence of absence of the analyte cannot be verified. Assignment of <i>R</i> requires approval by the appropriate WSRC data validation coordinator. |
| U | Material analyzed for but not detected. Analytical result reported is less than the sample quantitation limit. |
| NJ | The analysis indicates the presence of an analyte that has been tentatively identified and the associated numerical value represents its approximate concentration. |

Note: These are only some of the qualifiers present in the database. All modifiers associated with the data are published in the result tables of EPD/EMS' quarterly groundwater monitoring reports, the official repository of the data.

"EPA" Qualifiers - EPA STORET Codes

| | |
|---------|---|
| (Blank) | Data not remarked |
| C | The result is calculated. |
| I | The result is less than the ssEQL, but equal to or greater than the MDL. Always reported with an associated EPA functional Guideline Code qualifier of <i>U</i> . |
| K | The actual concentration is known to be less than the reported result. |
| L | The actual concentration is known to be less than the reported result. |
| O | Sample received by laboratory, but the analysis was lost or not performed. |
| Q | Sample was held beyond normal holding time prior to analysis. |

Note: These are only some of the qualifiers present in the database. All modifiers associated with the data are published in the result tables of EPD/EMS' quarterly groundwater monitoring reports, the official repository of the data.

Table B. Assessment Monitoring Results for Groundwater Wells

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| <u>SRS Coord.</u> | <u>Lat/Longitude</u> | <u>Screen Zone Elevation</u> | <u>Top of Casing</u> | <u>Casing</u> | <u>Pump</u> | <u>Screen Zone</u> |
|------------------------|----------------------------------|------------------------------|----------------------|---------------|-------------|--------------------|
| N 86079.6 E 45555.3 | 33.29048 Deg N -81.7141 Deg W | 162.1 - 141.1 ft msl | 192.4 ft msl | 4 " PVC | S | Upper |

SAMPLE DATE 03/08/00 10/02/00

FIELD DATA

| <u>Parameter</u> | <u>1st Half</u> | <u>2nd Half</u> | <u>Unit</u> |
|---------------------|-----------------|-----------------|-------------|
| Water Elevation | 161.9 | 161.3 | ft msl |
| pH | 5.2 | 5.1 | pH |
| Sp. Conductance | 22 | 19 | uS/cm |
| Water temperature | 20.4 | 20.8 | deg. C |
| Alkalinity as CaCO3 | 5 | 1 | mg/L |
| Turbidity | 1.3 | .7 | NTU |
| Volumes purged | 2.81 | 3.05 | well volume |
| Sampling code | | | |

ANALYTICAL DATA

Groundwater Protection Standard

| <u>SI</u> | <u>Parameter</u> | <u>1st Half</u> | <u>CLP</u> | <u>EPA</u> | <u>Filt.</u> | <u>2nd Half</u> | <u>CLP</u> | <u>EPA</u> | <u>Filt.</u> | <u>Unit</u> |
|-----------|--------------------------------|-----------------|------------|------------|--------------|-----------------|------------|------------|--------------|-------------|
| | 1,1,1,2-Tetrachloroethane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,1,1-Trichloroethane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,1,2,2-Tetrachloroethane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,1,2-Trichloroethane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,1-Dichloroethane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,1-Dichloroethylene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,1-Dichloropropene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,2,3-Trichloropropane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,2,4,5-Tetrachlorobenzene | | | | | | | | | |
| | 1,2,4-Trichlorobenzene | | | | | | | | | |
| | 1,2-Dibromo-3-chloropropane | <10 | U | | < EQL | <10 | U | | < EQL | ug/L |
| | 1,2-Dibromoethane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,2-Dichlorobenzene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,2-Dichloroethane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,2-Dichloropropane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,3,5-Trinitrobenzene | | | | | | | | | |
| | 1,3-Dichlorobenzene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,3-Dichloropropane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,3-Dinitrobenzene | | | | | | | | | |
| | 1,4-Dichlorobenzene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,4-Naphthoquinone | | | | | | | | | |
| | 1-Naphthylamine | | | | | | | | | |
| | 2,2-Dichloropropane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 2,2-Oxybis(1-chloropropane) | | | | | | | | | |
| | 2,3,4,6-Tetrachlorophenol | | | | | | | | | |
| | 2,4,5-T | | | | | | | | | |
| | 2,4,5-TP (Silvex) | | | | | | | | | |
| | 2,4,5-Trichlorophenol | | | | | | | | | |
| | 2,4,6-Trichlorophenol | | | | | | | | | |
| | 2,4-Dichlorophenol | | | | | | | | | |
| | 2,4-Dichlorophenoxyacetic acid | | | | | | | | | |
| | 2,4-Dimethyl phenol | | | | | | | | | |
| | 2,4-Dinitrophenol | | | | | | | | | |
| | 2,4-Dinitrotoluene | | | | | | | | | |
| | 2,6-Dichlorophenol | | | | | | | | | |
| | 2,6-Dinitrotoluene | | | | | | | | | |
| | 2-Acetylaminofluorene | | | | | | | | | |
| | 2-Chloronaphthalene | | | | | | | | | |
| | 2-Chlorophenol | | | | | | | | | |
| | 2-Hexanone | <20 | U | | < EQL | <20 | U | | < EQL | ug/L |
| | 2-Methyl-4,6-dinitrophenol | | | | | | | | | |
| | 2-Methylnaphthalene | | | | | | | | | |
| | 2-Naphthylamine | | | | | | | | | |
| | 2-Nitrophenol | | | | | | | | | |
| | 2-sec-Butyl-4,6-dinitrophenol | | | | | | | | | |
| | 3,3'-Dichlorobenzidine | | | | | | | | | |
| | 3,3'-Dimethylbenzidine | | | | | | | | | |
| | 3-Methylcholanthrene | | | | | | | | | |
| | 4-Aminobiphenyl | | | | | | | | | |
| | 4-Bromophenyl phenyl ether | | | | | | | | | |
| | 4-Chloro-m-cresol | | | | | | | | | |

Notes: Concentrations in bold italics exceed the groundwater protection standards listed in Appendix A. Bold italics were not assigned if the result was qualified with a "J", "R", "L", or "U" qualifier.

+ = exceeded the Groundwater Protection Standards listed in Appendix A.

Table B. Assessment Monitoring Results for Groundwater Wells (Cont.)

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ANALYTICAL DATA

Groundwater Protection Standard

| ST | Parameter | 1st Half | CLP | EPA | Filt. | 2nd Half | CLP | EPA | Filt. | Unit |
|----|--------------------------------------|----------|-----|-----|--------|----------|-----|-----|--------|------|
| | 4-Chloroaniline | | | | | | | | | |
| | 4-Chlorophenyl phenyl ether | | | | | | | | | |
| | 4-Nitrophenol | | | | | | | | | |
| | 5-Nitro-o-toluidine | | | | | | | | | |
| | 7,12-Dimethylbenz(a)anthracene | | | | | | | | | |
| | Acenaphthene | | | | | | | | | |
| | Acenaphthylene | | | | | | | | | |
| | Acetone | <20 | U | | < EQL | <20 | U | | < EQL | ug/L |
| | Acetonitrile (Methyl cyanide) | <200 | U | | < EQL | <200 | U | | < EQL | ug/L |
| | Acetophenone | | | | | | | | | |
| | Acrolein | <50 | U | | < EQL | <50 | U | | < EQL | ug/L |
| | Acrylonitrile | <10 | U | | < EQL | <10 | U | | < EQL | ug/L |
| | Aldrin | | | | | | | | | |
| | Allyl chloride | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Anthracene | | | | | | | | | |
| | Antimony, total recoverable | <100 | U | | < EQL | <100 | U | | < EQL | ug/L |
| | Arsenic, total recoverable | <10 | U | | < EQL | <3.21 | JU | | < EQL | ug/L |
| | Barium, total recoverable | 15.1 | | | < 2000 | 10.7 | | | < 2000 | ug/L |
| | Benzene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Benzo(a)anthracene | | | | | | | | | |
| | Benzo(a)pyrene | | | | | | | | | |
| | Benzo(b)fluoranthene | | | | | | | | | |
| | Benzo(g,h,i)perylene | | | | | | | | | |
| | Benzo(k)fluoranthene | | | | | | | | | |
| | Benzyl alcohol | | | | | | | | | |
| | Beryllium, total recoverable | .94 | J | I | NDD | <.13 | JU | | < EQL | ug/L |
| | Bis(2-chloroethoxy) methane | | | | | | | | | |
| | Bis(2-chloroethyl) ether | | | | | | | | | |
| | Bis(2-ethylhexyl) phthalate | | | | | | | | | |
| | Bromochloromethane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Bromodichloromethane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Bromoform | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Bromomethane (Methyl bromide) | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Butylbenzyl phthalate | | | | | | | | | |
| | Cadmium, total recoverable | <10 | U | | < EQL | <10 | U | | < EQL | ug/L |
| | Carbon disulfide | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Carbon tetrachloride | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Chlorobenzene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Chlorobenzilate | | | | | | | | | |
| | Chloroethane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Chloroethene (Vinyl chloride) | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Chloroform | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Chloromethane (Methyl chloride) | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Chloroprene | <20 | U | | < EQL | <20 | U | | < EQL | ug/L |
| | Chromium, total recoverable | <10 | U | | < EQL | <10 | U | | < EQL | ug/L |
| | Chrysene | | | | | | | | | |
| | Cobalt, total recoverable | <20 | U | | < EQL | <20 | U | | < EQL | ug/L |
| | Copper, total recoverable | 16.9 | J | I | NDD | 11.1 | J | I | NDD | ug/L |
| | Cyanide | | | | | | | | | |
| | Di-n-butyl phthalate | <11 | JU | Q | < EQL | <9.8 | U | | < EQL | ug/L |
| | Di-n-octyl phthalate | | | | | | | | | |
| | Diallate | | | | | | | | | |
| | Dibenz(a,h)anthracene | | | | | | | | | |
| | Dibenzofuran | | | | | | | | | |
| | Dibromochloromethane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Dibromomethane (Methylene bromide) | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Dichlorodifluoromethane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Dichloromethane (Methylene chloride) | 4.6 | J | I | NDD | <1.4 | U | | < EQL | ug/L |
| | Dieldrin | | | | | | | | | |
| | Diethyl phthalate | | | | | | | | | |
| | Dimethoate | | | | | | | | | |
| | Dimethyl phthalate | | | | | | | | | |
| | Diphenylamine | | | | | | | | | |
| | Disulfoton | | | | | | | | | |
| | Endosulfan I | | | | | | | | | |
| | Endosulfan II | | | | | | | | | |
| | Endosulfan sulfate | | | | | | | | | |
| | Endrin | | | | | | | | | |

Notes: Concentrations in bold italics exceed the groundwater protection standards listed in Appendix A. Bold italics were not assigned if the result was qualified with a "J", "R", "L", or "U" qualifier.

+ = exceeded the Groundwater Protection Standards listed in Appendix A.

Table B. Assessment Monitoring Results for Groundwater Wells (Cont.)

WELL LFW 28

ANALYTICAL DATA

Groundwater Protection Standard

| ST | Parameter | 1st Half | CLP | EPA | Flt. | 2nd Half | CLP | EPA | Flt. | Unit |
|----|---------------------------------|----------|-----|-----|-------|----------|-----|-----|-------|-------|
| | Endrin aldehyde | | | | | | | | | |
| | Ethyl methacrylate | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Ethyl methanesulfonate | | | | | | | | | |
| | Ethylbenzene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Famphur | | | | | | | | | |
| | Fluoranthene | | | | | | | | | |
| | Fluorene | | | | | | | | | |
| | Heptachlor | | | | | | | | | |
| | Heptachlor epoxide | | | | | | | | | |
| | Hexachlorobenzene | | | | | | | | | |
| | Hexachlorobutadiene | | | | | | | | | |
| | Hexachlorocyclopentadiene | | | | | | | | | |
| | Hexachloroethane | | | | | | | | | |
| | Hexachloropropene | | | | | | | | | |
| | Indeno(1,2,3-c,d)pyrene | | | | | | | | | |
| | Iodomethane (Methyl iodide) | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Isobutyl alcohol | <500 | U | | < EQL | <500 | U | | < EQL | ug/L |
| | Isodrin | | | | | | | | | |
| | Isophorone | | | | | | | | | |
| | Isosafrole | | | | | | | | | |
| | Kepone | | | | | | | | | |
| | Lead, total recoverable | 3.39 | J | I | NDD | <10 | U | | < EQL | ug/L |
| | Lindane | | | | | | | | | |
| | Mercury, total recoverable | <.5 | U | | < EQL | <.5 | U | | < EQL | ug/L |
| | Methacrylonitrile | <200 | U | | < EQL | <200 | U | | < EQL | ug/L |
| | Methapyrilene | | | | | | | | | |
| | Methoxychlor | | | | | | | | | |
| | Methyl ethyl ketone | <20 | U | | < EQL | <20 | U | | < EQL | ug/L |
| | Methyl isobutyl ketone | <10 | U | | < EQL | <10 | U | | < EQL | ug/L |
| | Methyl methacrylate | <20 | U | | < EQL | <20 | U | | < EQL | ug/L |
| | Methyl methanesulfonate | | | | | | | | | |
| | N-Nitrosodi-n-butylamine | | | | | | | | | |
| | N-Nitrosodiethylamine | | | | | | | | | |
| | N-Nitrosodimethylamine | | | | | | | | | |
| | N-Nitrosodiphenylamine | | | | | | | | | |
| | N-Nitrosodipropylamine | | | | | | | | | |
| | N-Nitrosomethylethylamine | | | | | | | | | |
| | N-Nitrosopiperidine | | | | | | | | | |
| | N-Nitrosopyrrolidine | | | | | | | | | |
| | Naphthalene | | | | | | | | | |
| | Nickel, total recoverable | <50 | U | | < EQL | <50 | U | | < EQL | ug/L |
| | Nitrobenzene | | | | | | | | | |
| | O,O,O-Triethyl phosphorothioate | | | | | | | | | |
| | PCB | | | | | | | | | |
| | PCB 1016 | | | | | | | | | |
| | PCB 1232 | | | | | | | | | |
| | PCB 1242 | | | | | | | | | |
| | PCB 1248 | | | | | | | | | |
| | PCB 1254 | | | | | | | | | |
| | PCB 1260 | | | | | | | | | |
| | Parathion | | | | | | | | | |
| | Parathion methyl | | | | | | | | | |
| | Pentachlorobenzene | | | | | | | | | |
| | Pentachloronitrobenzene | | | | | | | | | |
| | Pentachlorophenol | | | | | | | | | |
| | Phenacetin | | | | | | | | | |
| | Phenanthrene | | | | | | | | | |
| | Phenol | | | | | | | | | |
| | Phorate | | | | | | | | | |
| | Pronamid | | | | | | | | | |
| | Propionitrile | <200 | U | | < EQL | <200 | U | | < EQL | ug/L |
| | Pyrene | | | | | | | | | |
| | Safrole | | | | | | | | | |
| | Selenium, total recoverable | <10 | U | | < EQL | <10 | U | | < EQL | ug/L |
| | Silver, total recoverable | <20 | U | | < EQL | <20 | U | | < EQL | ug/L |
| | Specific conductance | 19.4 | | | < 37 | 17.6 | | | < 37 | uS/cm |
| | Styrene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Sulfide | | | | | | | | | |

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+ = exceeded the Groundwater Protection Standards listed in Appendix A.

Table B. Assessment Monitoring Results for Groundwater Wells (Cont.)

WELL LFW 28

ANALYTICAL DATA

Groundwater Protection Standard

| ST | Parameter | 1st Half | CLP | EPA | Filt. | 2nd Half | CLP | EPA | Filt. | Unit |
|----|-----------------------------|----------|-----|-----|-------|----------|-----|-----|-------|------|
| | Tetrachloroethylene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Thallium, total recoverable | <10 | U | | < EQL | 6.81 | J | I | NDD | ug/L |
| | Thionazin | | | | | | | | | |
| | Tin, total recoverable | <200 | U | | < EQL | | | | | ug/L |
| | Toluene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Toxaphene | | | | | | | | | |
| | Trichloroethylene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Trichlorofluoromethane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Vanadium, total recoverable | <10 | U | | < EQL | <10 | U | | < EQL | ug/L |
| | Vinyl acetate | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Xylenes | <10 | U | | < EQL | <10 | U | | < EQL | ug/L |
| | Zinc, total recoverable | 14 | J | I | NDD | <20 | U | | < EQL | ug/L |
| | alpha-Benzene hexachloride | | | | | | | | | |
| | alpha-Chlordane | | | | | | | | | |
| | beta-Benzene hexachloride | <.1 | U | | < EQL | <.095 | U | | < EQL | ug/L |
| | cis-1,2-Dichloroethylene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | cis-1,3-Dichloropropene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | delta-Benzene hexachloride | | | | | | | | | |
| | gamma-Chlordane | | | | | | | | | |
| | m-Cresol (3-Methylphenol) | | | | | | | | | |
| | m-Nitroaniline | | | | | | | | | |
| | o-Cresol (2-Methylphenol) | | | | | | | | | |
| | o-Nitroaniline | | | | | | | | | |
| | o-Toluidine | | | | | | | | | |
| | p,p"-DDD | | | | | | | | | |
| | p,p"-DDE | | | | | | | | | |
| | p,p"-DDT | | | | | | | | | |
| | p-Cresol (4-Methylphenol) | | | | | | | | | |
| | p-Dimethylaminoazobenzene | | | | | | | | | |
| | p-Nitroaniline | | | | | | | | | |
| | p-Phenylenediamine | | | | | | | | | |
| | pH | 6.21 | J | Q | NDD | 5.25 | J | Q | NDD | pH |
| | trans-1,2-Dichloroethylene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | trans-1,3-Dichloropropene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | trans-1,4-Dichloro-2-butene | <20 | U | | < EQL | <20 | U | | < EQL | ug/L |

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+ = exceeded the Groundwater Protection Standards listed in Appendix A.

Table B. Assessment Monitoring Results for Groundwater Wells (Cont.)

WELL LFW 30

| | | | | | | |
|------------------------|----------------------------------|------------------------------|----------------------|---------------|-------------|--------------------|
| <u>SRS Coord.</u> | <u>Lat/Longitude</u> | <u>Screen Zone Elevation</u> | <u>Top of Casing</u> | <u>Casing</u> | <u>Pump</u> | <u>Screen Zone</u> |
| N 86318.4 E 45170.9 | 33.29038 Deg N -81.7155 Deg W | 162.7 - 141.7 ft msl | 210 ft msl | 4 " PVC | S | Upper |

| | | |
|--------------------|----------|----------|
| SAMPLE DATE | 03/07/00 | 08/17/00 |
|--------------------|----------|----------|

FIELD DATA

| | | | |
|---------------------|-----------------|-----------------|-------------|
| <u>Parameter</u> | <u>1st Half</u> | <u>2nd Half</u> | <u>Unit</u> |
| Water Elevation | 163.12 | 162.4 | ft msl |
| pH | 5.3 | 5.6 | pH |
| Sp. Conductance | 17 | 17 | uS/cm |
| Water temperature | 19.1 | 23.9 | deg. C |
| Alkalinity as CaCO3 | 1 | 1 | mg/L |
| Turbidity | .7 | .6 | NTU |
| Volumes purged | 3.74 | 3.87 | well volume |
| Sampling code | | | |

ANALYTICAL DATA

Groundwater Protection Standard

| <u>SI</u> | <u>Parameter</u> | <u>1st Half</u> | <u>CLP</u> | <u>EPA</u> | <u>Flt.</u> | <u>2nd Half</u> | <u>CLP</u> | <u>EPA</u> | <u>Flt.</u> | <u>Unit</u> |
|-----------|--------------------------------|-----------------|------------|------------|-------------|-----------------|------------|------------|-------------|-------------|
| | 1,1,1,2-Tetrachloroethane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,1,1-Trichloroethane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,1,2,2-Tetrachloroethane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,1,2-Trichloroethane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,1-Dichloroethane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,1-Dichloroethylene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,1-Dichloropropene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,2,3-Trichloropropane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,2,4,5-Tetrachlorobenzene | | | | | | | | | |
| | 1,2,4-Trichlorobenzene | | | | | | | | | |
| | 1,2-Dibromo-3-chloropropane | <10 | U | | < EQL | <10 | U | | < EQL | ug/L |
| | 1,2-Dibromoethane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,2-Dichlorobenzene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,2-Dichloroethane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,2-Dichloropropane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,3,5-Trinitrobenzene | | | | | | | | | |
| | 1,3-Dichlorobenzene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,3-Dichloropropane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,3-Dinitrobenzene | | | | | | | | | |
| | 1,4-Dichlorobenzene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,4-Naphthoquinone | | | | | | | | | |
| | 1-Naphthylamine | | | | | | | | | |
| | 2,2-Dichloropropane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 2,2-Oxybis(1-chloropropane) | | | | | | | | | |
| | 2,3,4,6-Tetrachlorophenol | | | | | | | | | |
| | 2,4,5-T | | | | | | | | | |
| | 2,4,5-TP (Silvex) | | | | | | | | | |
| | 2,4,5-Trichlorophenol | | | | | | | | | |
| | 2,4,6-Trichlorophenol | | | | | | | | | |
| | 2,4-Dichlorophenol | | | | | | | | | |
| | 2,4-Dichlorophenoxyacetic acid | | | | | | | | | |
| | 2,4-Dimethyl phenol | | | | | | | | | |
| | 2,4-Dinitrophenol | | | | | | | | | |
| | 2,4-Dinitrotoluene | | | | | | | | | |
| | 2,6-Dichlorophenol | | | | | | | | | |
| | 2,6-Dinitrotoluene | | | | | | | | | |
| | 2-Acetylaminofluorene | | | | | | | | | |
| | 2-Chloronaphthalene | | | | | | | | | |
| | 2-Chlorophenol | | | | | | | | | |
| | 2-Hexanone | <20 | U | | < EQL | <20 | U | | < EQL | ug/L |
| | 2-Methyl-4,6-dinitrophenol | | | | | | | | | |
| | 2-Methylnaphthalene | | | | | | | | | |
| | 2-Naphthylamine | | | | | | | | | |
| | 2-Nitrophenol | | | | | | | | | |
| | 2-sec-Butyl-4,6-dinitrophenol | | | | | | | | | |
| | 3,3"-Dichlorobenzidine | | | | | | | | | |
| | 3,3"-Dimethylbenzidine | | | | | | | | | |
| | 3-Methylcholanthrene | | | | | | | | | |
| | 4-Aminobiphenyl | | | | | | | | | |
| | 4-Bromophenyl phenyl ether | | | | | | | | | |
| | 4-Chloro-m-cresol | | | | | | | | | |

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+ = exceeded the Groundwater Protection Standards listed in Appendix A.

Table B. Assessment Monitoring Results for Groundwater Wells (Cont.)

WELL LFW 30

ANALYTICAL DATA

Groundwater Protection Standard

| ST | Parameter | 1st Half | CLP | EPA | Filt. | 2nd Half | CLP | EPA | Filt. | Unit |
|----|--------------------------------------|----------|-----|-----|-------|----------|-----|-----|-------|------|
| | 4-Chloroaniline | | | | | | | | | |
| | 4-Chlorophenyl phenyl ether | | | | | | | | | |
| | 4-Nitrophenol | | | | | | | | | |
| | 5-Nitro-o-toluidine | | | | | | | | | |
| | 7,12-Dimethylbenz(a)anthracene | | | | | | | | | |
| | Acenaphthene | | | | | | | | | |
| | Acenaphthylene | | | | | | | | | |
| | Acetone | <20 | U | | < EQL | <20 | U | | < EQL | ug/L |
| | Acetonitrile (Methyl cyanide) | <200 | U | | < EQL | <200 | U | | < EQL | ug/L |
| | Acetophenone | | | | | | | | | |
| | Acrolein | <50 | U | | < EQL | <50 | U | | < EQL | ug/L |
| | Acrylonitrile | <10 | U | | < EQL | <10 | U | | < EQL | ug/L |
| | Aldrin | | | | | | | | | |
| | Allyl chloride | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Anthracene | | | | | | | | | |
| | Antimony, total recoverable | <100 | U | | < EQL | <100 | U | | < EQL | ug/L |
| | Arsenic, total recoverable | <10 | U | | < EQL | <10 | U | | < EQL | ug/L |
| | Barium, total recoverable | <5.48 | U | V | < EQL | 5.46 | J | I | NDD | ug/L |
| | Benzene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Benzo(a)anthracene | | | | | | | | | |
| | Benzo(a)pyrene | | | | | | | | | |
| | Benzo(b)fluoranthene | | | | | | | | | |
| | Benzo(g,h,i)perylene | | | | | | | | | |
| | Benzo(k)fluoranthene | | | | | | | | | |
| | Benzyl alcohol | | | | | | | | | |
| | Beryllium, total recoverable | <1 | U | | < EQL | <1 | U | | < EQL | ug/L |
| | Bis(2-chloroethoxy) methane | | | | | | | | | |
| | Bis(2-chloroethyl) ether | | | | | | | | | |
| | Bis(2-ethylhexyl) phthalate | | | | | | | | | |
| | Bromochloromethane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Bromodichloromethane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Bromoform | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Bromomethane (Methyl bromide) | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Butylbenzyl phthalate | | | | | | | | | |
| | Cadmium, total recoverable | <10 | U | | < EQL | <10 | U | | < EQL | ug/L |
| | Carbon disulfide | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Carbon tetrachloride | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Chlorobenzene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Chlorobenzilate | | | | | | | | | |
| | Chloroethane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Chloroethene (Vinyl chloride) | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Chloroform | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Chloromethane (Methyl chloride) | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Chloroprene | <20 | U | | < EQL | <20 | U | | < EQL | ug/L |
| | Chromium, total recoverable | <3.83 | JU | | < EQL | <10 | U | | < EQL | ug/L |
| | Chrysene | | | | | | | | | |
| | Cobalt, total recoverable | <20 | U | | < EQL | <20 | U | | < EQL | ug/L |
| | Copper, total recoverable | 13.2 | J | I | NDD | <17.4 | U | V | < EQL | ug/L |
| | Cyanide | | | | | | | | | |
| | Di-n-butyl phthalate | <11 | JU | L | < EQL | <10 | JU | Q | < EQL | ug/L |
| | Di-n-octyl phthalate | | | | | | | | | |
| | Diallate | | | | | | | | | |
| | Dibenz(a,h)anthracene | | | | | | | | | |
| | Dibenzofuran | | | | | | | | | |
| | Dibromochloromethane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Dibromomethane (Methylene bromide) | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Dichlorodifluoromethane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Dichloromethane (Methylene chloride) | <1.4 | U | | < EQL | <10 | U | | < EQL | ug/L |
| | Dieldrin | | | | | | | | | |
| | Diethyl phthalate | | | | | | | | | |
| | Dimethoate | | | | | | | | | |
| | Dimethyl phthalate | | | | | | | | | |
| | Diphenylamine | | | | | | | | | |
| | Disulfoton | | | | | | | | | |
| | Endosulfan I | | | | | | | | | |
| | Endosulfan II | | | | | | | | | |
| | Endosulfan sulfate | | | | | | | | | |
| | Endrin | | | | | | | | | |

Notes: Concentrations in bold italics exceed the groundwater protection standards listed in Appendix A. Bold italics were not assigned if the result was qualified with a "J", "R", "L", or "U" qualifier.

+ = exceeded the Groundwater Protection Standards listed in Appendix A.

Table B. Assessment Monitoring Results for Groundwater Wells (Cont.)
WELL LPW 30

| ANALYTICAL DATA | | | | | | |
|---------------------------------|---------------------------------|----------|-----|-------|------|------|
| Groundwater Protection Standard | | | | | | |
| SI | Parameter | 1st Half | CLP | EPA | Flt. | Unit |
| | Endrin aldehyde | | | | | |
| | Ethyl methacrylate | <5 | U | < EQL | <5 | U |
| | Ethyl methanesulfonate | | | | | |
| | Ethylbenzene | <5 | U | < EQL | <5 | U |
| | Famphur | | | | | |
| | Fluoranthene | | | | | |
| | Fluorene | | | | | |
| | Hepachlor | | | | | |
| | Hepachlor epoxide | | | | | |
| | Hexachlorobenzene | | | | | |
| | Hexachlorobutadiene | | | | | |
| | Hexachlorocyclopentadiene | | | | | |
| | Hexachloroethane | | | | | |
| | Hexachloropropene | | | | | |
| | Indeno(1,2,3-c,d)pyrene | <5 | U | < EQL | <5 | U |
| | Iodomethane (Methyl Iodide) | | | | | |
| | Isobutyl alcohol | <500 | U | < EQL | <500 | U |
| | Isodrin | | | | | |
| | Isophorone | | | | | |
| | Isosafrole | | | | | |
| | Kepone | | | | | |
| | Lead, total recoverable | <2.5 | JU | < EQL | 16 | < 18 |
| | Lindane | | | | | |
| | Mercury, total recoverable | <.5 | U | < EQL | <.5 | U |
| | Methacrylonitrile | <200 | U | < EQL | <200 | U |
| | Methapyrene | | | | | |
| | Methoxychlor | | | | | |
| | Methyl ethyl ketone | <20 | U | < EQL | <20 | U |
| | Methyl isobutyl ketone | <10 | U | < EQL | <10 | U |
| | Methyl methacrylate | <20 | U | < EQL | <20 | U |
| | Methyl methanesulfonate | | | | | |
| | N-Nitrosodi-n-butylamine | | | | | |
| | N-Nitrosodimethylamine | | | | | |
| | N-Nitrosodiphenylamine | | | | | |
| | N-Nitrosodipropylamine | | | | | |
| | N-Nitrosomethylethylamine | | | | | |
| | N-Nitrosopiperidine | | | | | |
| | N-Nitrosopyrrolidine | | | | | |
| | Naphthalene | | | | | |
| | Nickel, total recoverable | <50 | U | < EQL | <50 | U |
| | Nitrobenzene | | | | | |
| | O,O,O-Triethyl phosphorothioate | | | | | |
| | PCB | | | | | |
| | PCB 1016 | | | | | |
| | PCB 1232 | | | | | |
| | PCB 1242 | | | | | |
| | PCB 1248 | | | | | |
| | PCB 1254 | | | | | |
| | PCB 1260 | | | | | |
| | Parathion | | | | | |
| | Parathion methyl | | | | | |
| | Pentachlorobenzene | | | | | |
| | Pentachloronitrobenzene | | | | | |
| | Pentachlorophenol | | | | | |
| | Phenaceln | | | | | |
| | Phenanthrene | | | | | |
| | Phenol | | | | | |
| | Phorate | | | | | |
| | Pronamid | | | | | |
| | Propionitrile | <200 | U | < EQL | <200 | U |
| | Pyrene | | | | | |
| | Safrole | | | | | |
| | Selenium, total recoverable | <10 | U | < EQL | <10 | U |
| | Silver, total recoverable | <3.57 | JU | < EQL | <20 | U |
| | Specific conductance | 15.7 | | < 37 | 15.4 | |
| | Styrene | <5 | U | < EQL | <5 | U |
| | Sulfide | | | | | |

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+ = exceeded the Groundwater Protection Standards listed in Appendix A.

Table B. Assessment Monitoring Results for Groundwater Wells (Cont.)

WELL LFW 30

ANALYTICAL DATA

Groundwater Protection Standard

| ST | Parameter | 1st Half | CLP | EPA | Filt. | 2nd Half | CLP | EPA | Filt. | Unit |
|----|-----------------------------|----------|-----|-----|-------|----------|-----|-----|--------|------|
| | Tetrachloroethylene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Thallium, total recoverable | <6.28 | JU | V | < EQL | <10 | U | | < EQL | ug/L |
| | Thionazin | | | | | | | | | |
| | Tin, total recoverable | | | | | | | | | |
| | Toluene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Toxaphene | | | | | | | | | |
| | Trichloroethylene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Trichlorofluoromethane | 2.4 | J | I | NDD | <5 | U | | < EQL | ug/L |
| | Vanadium, total recoverable | <10 | U | | < EQL | <10 | U | | < EQL | ug/L |
| | Vinyl acetate | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Xylenes | <10 | U | | < EQL | <10 | U | | < EQL | ug/L |
| | Zinc, total recoverable | 18.4 | J | I | NDD | 24.6 | | | < 29.3 | ug/L |
| | alpha-Benzene hexachloride | | | | | | | | | |
| | alpha-Chlordane | | | | | | | | | |
| | beta-Benzene hexachloride | <.1 | U | | < EQL | <.095 | JU | Q | < EQL | ug/L |
| | cis-1,2-Dichloroethylene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | cis-1,3-Dichloropropene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | delta-Benzene hexachloride | | | | | | | | | |
| | gamma-Chlordane | | | | | | | | | |
| | m-Cresol (3-Methylphenol) | | | | | | | | | |
| | m-Nitroaniline | | | | | | | | | |
| | o-Cresol (2-Methylphenol) | | | | | | | | | |
| | o-Nitroaniline | | | | | | | | | |
| | o-Toluidine | | | | | | | | | |
| | p,p"-DDD | | | | | | | | | |
| | p,p"-DDE | | | | | | | | | |
| | p,p"-DDT | | | | | | | | | |
| | p-Cresol (4-Methylphenol) | | | | | | | | | |
| | p-Dimethylaminoazobenzene | | | | | | | | | |
| | p-Nitroaniline | | | | | | | | | |
| | p-Phenylenediamine | | | | | | | | | |
| | pH | 5.64 | J | Q | NDD | 5.18 | J | Q | NDD | pH |
| | trans-1,2-Dichloroethylene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | trans-1,3-Dichloropropene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | trans-1,4-Dichloro-2-butene | <20 | U | | < EQL | <20 | U | | < EQL | ug/L |

Notes: Concentrations in bold italics exceed the groundwater protection standards listed in Appendix A. Bold italics were not assigned if the result was qualified with a "J", "R", "L", or "U" qualifier.

+ = exceeded the Groundwater Protection Standards listed in Appendix A.

Table B. Assessment Monitoring Results for Groundwater Wells (Cont.)

WELL LFW 31

| SRS Coord. | Lat/Longitude | Screen Zone Elevation | Top of Casing | Casing | Pump | Screen Zone |
|----------------------|----------------------------------|-----------------------|---------------|---------|------|-------------|
| N 86262.2 E 44869 | 33.28976 Deg N -81.7162 Deg W | 166 - 145 ft msl | 229.3 ft msl | 4 " PVC | S | Upper |

SAMPLE DATE 03/07/00 08/17/00

FIELD DATA

| Parameter | 1st Half | 2nd Half | Unit |
|---------------------|----------|----------|-------------|
| Water Elevation | 162.94 | 162.19 | ft msl |
| pH | 5.2 | 5.6 | pH |
| Sp. Conductance | 15 | 19 | uS/cm |
| Water temperature | 18.5 | 22.4 | deg. C |
| Alkalinity as CaCO3 | 2 | 2 | mg/L |
| Turbidity | 1.7 | .6 | NTU |
| Volumes purged | 4.15 | 2.53 | well volume |
| Sampling code | | | |

ANALYTICAL DATA

Groundwater Protection Standard

| ST | Parameter | 1st Half | CLP | EPA | Filt. | 2nd Half | CLP | EPA | Filt. | Unit |
|----|--------------------------------|----------|-----|-----|-------|----------|-----|-----|-------|------|
| | 1,1,1,2-Tetrachloroethane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,1,1-Trichloroethane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,1,2,2-Tetrachloroethane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,1,2-Trichloroethane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,1-Dichloroethane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,1-Dichloroethylene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,1-Dichloropropene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,2,3-Trichloropropane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,2,4,5-Tetrachlorobenzene | | | | | | | | | |
| | 1,2,4-Trichlorobenzene | | | | | | | | | |
| | 1,2-Dibromo-3-chloropropane | <10 | U | | < EQL | <10 | U | | < EQL | ug/L |
| | 1,2-Dibromoethane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,2-Dichlorobenzene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,2-Dichloroethane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,2-Dichloropropane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,3,5-Trinitrobenzene | | | | | | | | | |
| | 1,3-Dichlorobenzene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,3-Dichloropropane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,3-Dinitrobenzene | | | | | | | | | |
| | 1,4-Dichlorobenzene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,4-Naphthoquinone | | | | | | | | | |
| | 1-Naphthylamine | | | | | | | | | |
| | 2,2-Dichloropropane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 2,2-Oxybis(1-chloropropane) | | | | | | | | | |
| | 2,3,4,6-Tetrachlorophenol | | | | | | | | | |
| | 2,4,5-T | | | | | | | | | |
| | 2,4,5-TP (Silvex) | | | | | | | | | |
| | 2,4,5-Trichlorophenol | | | | | | | | | |
| | 2,4,6-Trichlorophenol | | | | | | | | | |
| | 2,4-Dichlorophenol | | | | | | | | | |
| | 2,4-Dichlorophenoxyacetic acid | | | | | | | | | |
| | 2,4-Dimethyl phenol | | | | | | | | | |
| | 2,4-Dinitrophenol | | | | | | | | | |
| | 2,4-Dinitrotoluene | | | | | | | | | |
| | 2,6-Dichlorophenol | | | | | | | | | |
| | 2,6-Dinitrotoluene | | | | | | | | | |
| | 2-Acetylaminofluorene | | | | | | | | | |
| | 2-Chloronaphthalene | | | | | | | | | |
| | 2-Chlorophenol | | | | | | | | | |
| | 2-Hexanone | <20 | U | | < EQL | <20 | U | | < EQL | ug/L |
| | 2-Methyl-4,6-dinitrophenol | | | | | | | | | |
| | 2-Methylnaphthalene | | | | | | | | | |
| | 2-Naphthylamine | | | | | | | | | |
| | 2-Nitrophenol | | | | | | | | | |
| | 2-sec-Butyl-4,6-dinitrophenol | | | | | | | | | |
| | 3,3"-Dichlorobenzidine | | | | | | | | | |
| | 3,3"-Dimethylbenzidine | | | | | | | | | |
| | 3-Methylcholanthrene | | | | | | | | | |
| | 4-Aminobiphenyl | | | | | | | | | |
| | 4-Bromophenyl phenyl ether | | | | | | | | | |
| | 4-Chloro-m-cresol | | | | | | | | | |

Notes: Concentrations in bold italics exceed the groundwater protection standards listed in Appendix A. Bold italics were not assigned if the result was qualified with a "J", "R", "L", or "U" qualifier.

+ = exceeded the Groundwater Protection Standards listed in Appendix A.

Table B. Assessment Monitoring Results for Groundwater Wells (Cont.)

WELL LFW 31

ANALYTICAL DATA

Groundwater Protection Standard

| ST | Parameter | 1st Half | CLP | EPA | Filt. | 2nd Half | CLP | EPA | Filt. | Unit |
|----|--------------------------------------|----------|-----|-----|-------|----------|-----|-----|--------|------|
| | 4-Chloroaniline | | | | | | | | | |
| | 4-Chlorophenyl phenyl ether | | | | | | | | | |
| | 4-Nitrophenol | | | | | | | | | |
| | 5-Nitro-o-toluidine | | | | | | | | | |
| | 7,12-Dimethylbenz(a)anthracene | | | | | | | | | |
| | Acenaphthene | | | | | | | | | |
| | Acenaphthylene | | | | | | | | | |
| | Acetone | <20 | U | | < EQL | <20 | U | | < EQL | ug/L |
| | Acetonitrile (Methyl cyanide) | <200 | U | | < EQL | <200 | U | | < EQL | ug/L |
| | Acetophenone | | | | | | | | | |
| | Acrolein | <50 | U | | < EQL | <50 | U | | < EQL | ug/L |
| | Acrylonitrile | <10 | U | | < EQL | <10 | U | | < EQL | ug/L |
| | Aldrin | | | | | | | | | |
| | Allyl chloride | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Anthracene | | | | | | | | | |
| | Antimony, total recoverable | <100 | U | | < EQL | <100 | U | | < EQL | ug/L |
| | Arsenic, total recoverable | <10 | U | | < EQL | <10 | U | | < EQL | ug/L |
| | Barium, total recoverable | 4.29 | J | I | NDD | 5.09 | J | I | NDD | ug/L |
| | Benzene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Benzo(a)anthracene | | | | | | | | | |
| | Benzo(a)pyrene | | | | | | | | | |
| | Benzo(b)fluoranthene | | | | | | | | | |
| | Benzo(g,h,i)perylene | | | | | | | | | |
| | Benzo(k)fluoranthene | | | | | | | | | |
| | Benzyl alcohol | | | | | | | | | |
| | Beryllium, total recoverable | <1 | U | | < EQL | <1 | U | | < EQL | ug/L |
| | Bis(2-chloroethoxy) methane | | | | | | | | | |
| | Bis(2-chloroethyl) ether | | | | | | | | | |
| | Bis(2-ethylhexyl) phthalate | | | | | | | | | |
| | Bromochloromethane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Bromodichloromethane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Bromoform | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Bromomethane (Methyl bromide) | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Butylbenzyl phthalate | | | | | | | | | |
| | Cadmium, total recoverable | <10 | U | | < EQL | <10 | U | | < EQL | ug/L |
| | Carbon disulfide | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Carbon tetrachloride | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Chlorobenzene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Chlorobenzilate | | | | | | | | | |
| | Chloroethane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Chloroethene (Vinyl chloride) | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Chloroform | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Chloromethane (Methyl chloride) | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Chloroprene | <20 | U | | < EQL | <20 | U | | < EQL | ug/L |
| | Chromium, total recoverable | <10 | U | | < EQL | <10 | U | | < EQL | ug/L |
| | Chrysene | | | | | | | | | |
| | Cobalt, total recoverable | <20 | U | | < EQL | <20 | U | | < EQL | ug/L |
| + | Copper, total recoverable | 6.78 | J | I | NDD | 40.8 | | | > 29.8 | ug/L |
| | Cyanide | | | | | | | | | |
| | Di-n-butyl phthalate | <10 | JU | L | < EQL | <9.5 | JU | Q | < EQL | ug/L |
| | Di-n-octyl phthalate | | | | | | | | | |
| | Diallate | | | | | | | | | |
| | Dibenz(a,h)anthracene | | | | | | | | | |
| | Dibenzofuran | | | | | | | | | |
| | Dibromochloromethane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Dibromomethane (Methylene bromide) | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Dichlorodifluoromethane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Dichloromethane (Methylene chloride) | <2.5 | U | V | < EQL | <10 | U | | < EQL | ug/L |
| | Dieldrin | | | | | | | | | |
| | Diethyl phthalate | | | | | | | | | |
| | Dimethoate | | | | | | | | | |
| | Dimethyl phthalate | | | | | | | | | |
| | Diphenylamine | | | | | | | | | |
| | Disulfoton | | | | | | | | | |
| | Endosulfan I | | | | | | | | | |
| | Endosulfan II | | | | | | | | | |
| | Endosulfan sulfate | | | | | | | | | |
| | Endrin | | | | | | | | | |

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+ = exceeded the Groundwater Protection Standards listed in Appendix A.

Table B. Assessment Monitoring Results for Groundwater Wells (Cont.)

WELL LFW 31

ANALYTICAL DATA

Groundwater Protection Standard

| ST | Parameter | 1st Half | CLP | EPA | Filt. | 2nd Half | CLP | EPA | Filt. | Unit |
|----|---------------------------------|----------|-----|-----|-------|----------|-----|-----|-------|-------|
| | Endrin aldehyde | | | | | | | | | |
| | Ethyl methacrylate | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Ethyl methanesulfonate | | | | | | | | | |
| | Ethylbenzene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Famphur | | | | | | | | | |
| | Fluoranthene | | | | | | | | | |
| | Fluorene | | | | | | | | | |
| | Heptachlor | | | | | | | | | |
| | Heptachlor epoxide | | | | | | | | | |
| | Hexachlorobenzene | | | | | | | | | |
| | Hexachlorobutadiene | | | | | | | | | |
| | Hexachlorocyclopentadiene | | | | | | | | | |
| | Hexachloroethane | | | | | | | | | |
| | Hexachloropropene | | | | | | | | | |
| | Indeno(1,2,3-c,d)pyrene | | | | | | | | | |
| | Iodomethane (Methyl iodide) | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Isobutyl alcohol | <500 | U | | < EQL | <500 | U | | < EQL | ug/L |
| | Isodrin | | | | | | | | | |
| | Isophorone | | | | | | | | | |
| | Isosafrole | | | | | | | | | |
| | Kepone | | | | | | | | | |
| | Lead, total recoverable | <3.05 | JU | | < EQL | <10 | U | | < EQL | ug/L |
| | Lindane | | | | | | | | | |
| | Mercury, total recoverable | <.5 | U | | < EQL | <.5 | U | | < EQL | ug/L |
| | Methacrylonitrile | <200 | U | | < EQL | <200 | U | | < EQL | ug/L |
| | Methapyrilene | | | | | | | | | |
| | Methoxychlor | | | | | | | | | |
| | Methyl ethyl ketone | <20 | U | | < EQL | <20 | U | | < EQL | ug/L |
| | Methyl isobutyl ketone | <10 | U | | < EQL | <10 | U | | < EQL | ug/L |
| | Methyl methacrylate | <20 | U | | < EQL | <20 | U | | < EQL | ug/L |
| | Methyl methanesulfonate | | | | | | | | | |
| | N-Nitrosodi-n-butylamine | | | | | | | | | |
| | N-Nitrosodiethylamine | | | | | | | | | |
| | N-Nitrosodimethylamine | | | | | | | | | |
| | N-Nitrosodiphenylamine | | | | | | | | | |
| | N-Nitrosodipropylamine | | | | | | | | | |
| | N-Nitrosomethylethylamine | | | | | | | | | |
| | N-Nitrosopiperidine | | | | | | | | | |
| | N-Nitrosopyrrolidine | | | | | | | | | |
| | Naphthalene | | | | | | | | | |
| | Nickel, total recoverable | <50 | U | | < EQL | <50 | U | | < EQL | ug/L |
| | Nitrobenzene | | | | | | | | | |
| | O,O,O-Triethyl phosphorothioate | | | | | | | | | |
| | PCB | | | | | | | | | |
| | PCB 1016 | | | | | | | | | |
| | PCB 1232 | | | | | | | | | |
| | PCB 1242 | | | | | | | | | |
| | PCB 1248 | | | | | | | | | |
| | PCB 1254 | | | | | | | | | |
| | PCB 1260 | | | | | | | | | |
| | Parathion | | | | | | | | | |
| | Parathion methyl | | | | | | | | | |
| | Pentachlorobenzene | | | | | | | | | |
| | Pentachloronitrobenzene | | | | | | | | | |
| | Pentachlorophenol | | | | | | | | | |
| | Phenacetin | | | | | | | | | |
| | Phenanthrene | | | | | | | | | |
| | Phenol | | | | | | | | | |
| | Phorate | | | | | | | | | |
| | Pronamid | | | | | | | | | |
| | Propionitrile | <200 | U | | < EQL | <200 | U | | < EQL | ug/L |
| | Pyrene | | | | | | | | | |
| | Safrole | | | | | | | | | |
| | Selenium, total recoverable | 12.3 | | | < 50 | <10 | U | | < EQL | ug/L |
| | Silver, total recoverable | 46.1 | | | < 100 | <20 | U | | < EQL | ug/L |
| | Specific conductance | 14.8 | | | < 37 | 18.9 | | | < 37 | uS/cm |
| | Styrene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Sulfide | | | | | | | | | |

Notes: Concentrations in bold italics exceed the groundwater protection standards listed in Appendix A. Bold italics were not assigned if the result was qualified with a "J", "R", "L", or "U" qualifier.

+ = exceeded the Groundwater Protection Standards listed in Appendix A.

Table B. Assessment Monitoring Results for Groundwater Wells (Cont.)

WELL LFW 31

ANALYTICAL DATA

Groundwater Protection Standard

| ST | Parameter | 1st Half | CLP | EPA | Filt. | 2nd Half | CLP | EPA | Filt. | Unit |
|----|-----------------------------|----------|-----|-----|-------|-------------|-----|-----|--------|------|
| | Tetrachloroethylene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Thallium, total recoverable | <5.65 | JU | V | < EQL | <10 | U | | < EQL | ug/L |
| | Thionazin | | | | | | | | | |
| | Tin, total recoverable | | | | | | | | | |
| | Toluene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Toxaphene | | | | | | | | | |
| | Trichloroethylene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Trichlorofluoromethane | 1.5 | J | IL | NDD | 3.7 | J | I | NDD | ug/L |
| | Vanadium, total recoverable | <10 | U | | < EQL | <10 | U | | < EQL | ug/L |
| | Vinyl acetate | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Xylenes | <10 | U | | < EQL | <10 | U | | < EQL | ug/L |
| + | Zinc, total recoverable | 11 | J | I | NDD | 40.2 | | | > 29.3 | ug/L |
| | alpha-Benzene hexachloride | | | | | | | | | |
| | alpha-Chlordane | | | | | | | | | |
| | beta-Benzene hexachloride | <.1 | U | | < EQL | <.095 | JU | Q | < EQL | ug/L |
| | cis-1,2-Dichloroethylene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | cis-1,3-Dichloropropene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | delta-Benzene hexachloride | | | | | | | | | |
| | gamma-Chlordane | | | | | | | | | |
| | m-Cresol (3-Methylphenol) | | | | | | | | | |
| | m-Nitroaniline | | | | | | | | | |
| | o-Cresol (2-Methylphenol) | | | | | | | | | |
| | o-Nitroaniline | | | | | | | | | |
| | o-Toluidine | | | | | | | | | |
| | p,p'-DDD | | | | | | | | | |
| | p,p'-DDE | | | | | | | | | |
| | p,p'-DDT | | | | | | | | | |
| | p-Cresol (4-Methylphenol) | | | | | | | | | |
| | p-Dimethylaminoazobenzene | | | | | | | | | |
| | p-Nitroaniline | | | | | | | | | |
| | p-Phenylenediamine | | | | | | | | | |
| | pH | 5.73 | J | Q | NDD | 5.24 | J | Q | NDD | pH |
| | trans-1,2-Dichloroethylene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | trans-1,3-Dichloropropene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | trans-1,4-Dichloro-2-butene | <20 | U | | < EQL | <20 | U | | < EQL | ug/L |

Notes: Concentrations in bold italics exceed the groundwater protection standards listed in Appendix A. Bold italics were not assigned if the result was qualified with a "J", "R", "L", or "U" qualifier.

+ = exceeded the Groundwater Protection Standards listed in Appendix A.

Table B. Assessment Monitoring Results for Groundwater Wells (Cont.)

WELL LFW 32

| SRS Coord. | Lat/Longitude | Screen Zone Elevation | Top of Casing | Casing | Pump | Screen Zone |
|------------------------|----------------------------------|-----------------------|---------------|---------|------|-------------|
| N 85836.8 E 44935.9 | 33.28893 Deg N -81.7152 Deg W | 165.3 - 144.3 ft msl | 223.7 ft msl | 4 " PVC | S | Upper |

SAMPLE DATE 03/08/00 10/02/00

FIELD DATA

| Parameter | 1st Half | 2nd Half | Unit |
|---------------------|----------|----------|-------------|
| Water Elevation | 160.99 | 160.45 | ft msl |
| pH | 4.2 | 4.8 | pH |
| Sp. Conductance | 27 | 29 | uS/cm |
| Water temperature | 19.4 | 20.7 | deg. C |
| Alkalinity as CaCO3 | 0 | 0 | mg/L |
| Turbidity | .9 | .6 | NTU |
| Volumes purged | 3.51 | 3.25 | well volume |
| Sampling code | | | |

ANALYTICAL DATA

Groundwater Protection Standard

| SI | Parameter | 1st Half | CLP | EPA | Filt. | 2nd Half | CLP | EPA | Filt. | Unit |
|----|--------------------------------|----------|-----|-----|-------|----------|-----|-----|-------|------|
| | 1,1,1,2-Tetrachloroethane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,1,1-Trichloroethane | 4.6 | J | I | NDD | 6.3 | J | K | NDD | ug/L |
| | 1,1,2,2-Tetrachloroethane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,1,2-Trichloroethane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,1-Dichloroethane | 2.5 | J | I | NDD | 4.6 | J | IK | NDD | ug/L |
| | 1,1-Dichloroethylene | <5 | U | | < EQL | 1.1 | J | IK | NDD | ug/L |
| | 1,1-Dichloropropene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,2,3-Trichloropropane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,2,4,5-Tetrachlorobenzene | <53 | JU | Q | < EQL | | | | | ug/L |
| | 1,2,4-Trichlorobenzene | <11 | JU | Q | < EQL | | | | | ug/L |
| | 1,2-Dibromo-3-chloropropane | <10 | U | | < EQL | <10 | U | | < EQL | ug/L |
| | 1,2-Dibromoethane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,2-Dichlorobenzene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,2-Dichloroethane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,2-Dichloropropane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,3,5-Trinitrobenzene | <10 | JU | Q | < EQL | | | | | ug/L |
| | 1,3-Dichlorobenzene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,3-Dichloropropane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,3-Dinitrobenzene | <50 | JU | Q | < EQL | | | | | ug/L |
| | 1,4-Dichlorobenzene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,4-Naphthoquinone | <11 | JU | Q | < EQL | | | | | ug/L |
| | 1-Naphthylamine | <11 | JU | Q | < EQL | | | | | ug/L |
| | 2,2-Dichloropropane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 2,2-Oxybis(1-chloropropane) | <11 | JU | Q | < EQL | | | | | ug/L |
| | 2,3,4,6-Tetrachlorophenol | <11 | JU | Q | < EQL | | | | | ug/L |
| | 2,4,5-T | <2 | U | Q | < EQL | | | | | ug/L |
| | 2,4,5-TP (Silvex) | <2 | U | Q | < EQL | | | | | ug/L |
| | 2,4,5-Trichlorophenol | <11 | JU | Q | < EQL | | | | | ug/L |
| | 2,4,6-Trichlorophenol | <26 | JU | Q | < EQL | | | | | ug/L |
| | 2,4-Dichlorophenol | <11 | JU | Q | < EQL | | | | | ug/L |
| | 2,4-Dichlorophenoxyacetic acid | <2 | U | Q | < EQL | | | | | ug/L |
| | 2,4-Dimethyl phenol | <11 | JU | Q | < EQL | | | | | ug/L |
| | 2,4-Dinitrophenol | <26 | JU | Q | < EQL | | | | | ug/L |
| | 2,4-Dinitrotoluene | <11 | JU | Q | < EQL | | | | | ug/L |
| | 2,6-Dichlorophenol | <25 | JU | Q | < EQL | | | | | ug/L |
| | 2,6-Dinitrotoluene | <11 | JU | Q | < EQL | | | | | ug/L |
| | 2-Acetylaminofluorene | <11 | JU | Q | < EQL | | | | | ug/L |
| | 2-Chloronaphthalene | <11 | JU | Q | < EQL | | | | | ug/L |
| | 2-Chlorophenol | <11 | JU | Q | < EQL | | | | | ug/L |
| | 2-Hexanone | <20 | U | | < EQL | <20 | U | | < EQL | ug/L |
| | 2-Methyl-4,6-dinitrophenol | <26 | JU | Q | < EQL | | | | | ug/L |
| | 2-Methylnaphthalene | <11 | JU | Q | < EQL | | | | | ug/L |
| | 2-Naphthylamine | <11 | JU | Q | < EQL | | | | | ug/L |
| | 2-Nitrophenol | <11 | JU | Q | < EQL | | | | | ug/L |
| | 2-sec-Butyl-4,6-dinitrophenol | <2 | U | Q | < EQL | | | | | ug/L |
| | 3,3"-Dichlorobenzidine | <11 | JU | Q | < EQL | | | | | ug/L |
| | 3,3"-Dimethylbenzidine | <21 | JU | Q | < EQL | | | | | ug/L |
| | 3-Methylcholanthrene | <11 | JU | Q | < EQL | | | | | ug/L |
| | 4-Aminobiphenyl | <11 | JU | Q | < EQL | | | | | ug/L |
| | 4-Bromophenyl phenyl ether | <11 | JU | Q | < EQL | | | | | ug/L |
| | 4-Chloro-m-cresol | <11 | JU | Q | < EQL | | | | | ug/L |

Notes: Concentrations in bold italics exceed the groundwater protection standards listed in Appendix A. Bold italics were not assigned if the result was qualified with a "J", "R", "L", or "U" qualifier.

+ = exceeded the Groundwater Protection Standards listed in Appendix A.

Table B. Assessment Monitoring Results for Groundwater Wells (Cont.)

WELL LFW 32

ANALYTICAL DATA

Groundwater Protection Standard

| ST | Parameter | 1st Half | CLP | EPA | Filt. | 2nd Half | CLP | EPA | Filt. | Unit |
|-----------|--------------------------------------|-----------------|------------|------------|--------------|-----------------|------------|------------|--------------|-------------|
| | 4-Chloroaniline | <11 | JU | Q | < EQL | | | | | ug/L |
| | 4-Chlorophenyl phenyl ether | <11 | JU | Q | < EQL | | | | | ug/L |
| | 4-Nitrophenol | <26 | JU | Q | < EQL | | | | | ug/L |
| | 5-Nitro-o-toluidine | <11 | JU | Q | < EQL | | | | | ug/L |
| | 7,12-Dimethylbenz(a)anthracene | <11 | JU | Q | < EQL | | | | | ug/L |
| | Acenaphthene | <11 | JU | Q | < EQL | | | | | ug/L |
| | Acenaphthylene | <11 | JU | Q | < EQL | | | | | ug/L |
| | Acetone | <20 | U | | < EQL | <20 | U | | < EQL | ug/L |
| | Acetonitrile (Methyl cyanide) | <200 | U | | < EQL | <200 | U | | < EQL | ug/L |
| | Acetophenone | <11 | JU | Q | < EQL | | | | | ug/L |
| | Acrolein | <50 | U | | < EQL | <50 | U | | < EQL | ug/L |
| | Acrylonitrile | <10 | U | | < EQL | <10 | U | | < EQL | ug/L |
| | Aldrin | <.1 | U | | < EQL | | | | | ug/L |
| | Allyl chloride | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Anthracene | <11 | JU | Q | < EQL | | | | | ug/L |
| | Antimony, total recoverable | <100 | U | | < EQL | <100 | U | | < EQL | ug/L |
| | Arsenic, total recoverable | <10 | U | | < EQL | <4.21 | JU | | < EQL | ug/L |
| | Barium, total recoverable | 10.9 | | | < 2000 | 11.4 | | | < 2000 | ug/L |
| | Benzene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Benzo(a)anthracene | <11 | JU | Q | < EQL | | | | | ug/L |
| | Benzo(a)pyrene | <11 | JU | Q | < EQL | | | | | ug/L |
| | Benzo(b)fluoranthene | <11 | JU | Q | < EQL | | | | | ug/L |
| | Benzo(g,h,i)perylene | <11 | JU | Q | < EQL | | | | | ug/L |
| | Benzo(k)fluoranthene | <11 | JU | Q | < EQL | | | | | ug/L |
| | Benzyl alcohol | | | | | | | | | |
| | Beryllium, total recoverable | <1 | U | | < EQL | <.19 | JU | | < EQL | ug/L |
| | Bis(2-chloroethoxy) methane | <11 | JU | Q | < EQL | | | | | ug/L |
| | Bis(2-chloroethyl) ether | <11 | JU | Q | < EQL | | | | | ug/L |
| | Bis(2-ethylhexyl) phthalate | <11 | JU | Q | < EQL | | | | | ug/L |
| | Bromochloromethane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Bromodichloromethane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Bromoform | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Bromomethane (Methyl bromide) | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Butylbenzyl phthalate | <11 | JU | Q | < EQL | | | | | ug/L |
| | Cadmium, total recoverable | <10 | U | | < EQL | <10 | U | | < EQL | ug/L |
| | Carbon disulfide | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Carbon tetrachloride | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Chlorobenzene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Chlorobenzilate | <11 | JU | Q | < EQL | | | | | ug/L |
| | Chloroethane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Chloroethene (Vinyl chloride) | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Chloroform | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Chloromethane (Methyl chloride) | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Chloroprene | <20 | U | | < EQL | <20 | U | | < EQL | ug/L |
| | Chromium, total recoverable | <10 | U | | < EQL | <10 | U | | < EQL | ug/L |
| | Chrysene | <11 | JU | Q | < EQL | | | | | ug/L |
| | Cobalt, total recoverable | <20 | U | | < EQL | <20 | U | | < EQL | ug/L |
| | Copper, total recoverable | 10.9 | J | I | NDD | <3.43 | JU | | < EQL | ug/L |
| | Cyanide | 10 | R | Q | Rejected | | | | | ug/L |
| | Di-n-butyl phthalate | <11 | JU | Q | < EQL | <9.5 | U | | < EQL | ug/L |
| | Di-n-octyl phthalate | <11 | JU | Q | < EQL | | | | | ug/L |
| | Diallate | <11 | JU | Q | < EQL | | | | | ug/L |
| | Dibenz(a,h)anthracene | <11 | JU | Q | < EQL | | | | | ug/L |
| | Dibenzofuran | <11 | JU | Q | < EQL | | | | | ug/L |
| | Dibromochloromethane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Dibromomethane (Methylene bromide) | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Dichlorodifluoromethane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Dichloromethane (Methylene chloride) | 3.7 | J | I | NDD | <4.7 | U | | < EQL | ug/L |
| | Dieldrin | <.21 | U | | < EQL | | | | | ug/L |
| | Diethyl phthalate | <11 | JU | Q | < EQL | | | | | ug/L |
| | Dimethoate | <11 | JU | Q | < EQL | | | | | ug/L |
| | Dimethyl phthalate | <11 | JU | Q | < EQL | | | | | ug/L |
| | Diphenylamine | <11 | JU | Q | < EQL | | | | | ug/L |
| | Disulfoton | <11 | JU | Q | < EQL | | | | | ug/L |
| | Endosulfan I | <.1 | U | | < EQL | | | | | ug/L |
| | Endosulfan II | <.21 | U | | < EQL | | | | | ug/L |
| | Endosulfan sulfate | <.21 | U | | < EQL | | | | | ug/L |
| | Endrin | <.21 | U | | < EQL | | | | | ug/L |

Notes: Concentrations in bold italics exceed the groundwater protection standards listed in Appendix A. Bold italics were not assigned if the result was qualified with a "J", "R", "L", or "U" qualifier.

+ = exceeded the Groundwater Protection Standards listed in Appendix A.

Table B. Assessment Monitoring Results for Groundwater Wells (Cont.)

WELL LFW 32

ANALYTICAL DATA

Groundwater Protection Standard

| ST | Parameter | 1st Half | CLP | EPA | Filt. | 2nd Half | CLP | EPA | Filt. | Unit |
|----|---------------------------------|----------|-----|-----|----------|----------|-----|-----|-------|-------|
| | Endrin aldehyde | <.21 | U | | < EQL | | | | | ug/L |
| | Ethyl methacrylate | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Ethyl methanesulfonate | <11 | JU | Q | < EQL | | | | | ug/L |
| | Ethylbenzene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Famphur | <200 | JU | Q | < EQL | | | | | ug/L |
| | Fluoranthene | <11 | JU | Q | < EQL | | | | | ug/L |
| | Fluorene | <11 | JU | Q | < EQL | | | | | ug/L |
| | Heptachlor | <.1 | U | | < EQL | | | | | ug/L |
| | Heptachlor epoxide | <.1 | U | | < EQL | | | | | ug/L |
| | Hexachlorobenzene | <11 | JU | Q | < EQL | | | | | ug/L |
| | Hexachlorobutadiene | <11 | JU | Q | < EQL | | | | | ug/L |
| | Hexachlorocyclopentadiene | <11 | JU | Q | < EQL | | | | | ug/L |
| | Hexachloroethane | <11 | JU | Q | < EQL | | | | | ug/L |
| | Hexachloropropene | <53 | JU | Q | < EQL | | | | | ug/L |
| | Indeno(1,2,3-c,d)pyrene | <11 | JU | Q | < EQL | | | | | ug/L |
| | Iodomethane (Methyl iodide) | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Isobutyl alcohol | <500 | U | | < EQL | <500 | U | | < EQL | ug/L |
| | Isodrin | <11 | JU | Q | < EQL | | | | | ug/L |
| | Isophorone | <11 | JU | Q | < EQL | | | | | ug/L |
| | Isosafrole | <11 | JU | Q | < EQL | | | | | ug/L |
| | Kepone | <11 | JU | Q | < EQL | | | | | ug/L |
| | Lead, total recoverable | <2.88 | JU | | < EQL | <10 | U | | < EQL | ug/L |
| | Lindane | <.1 | U | | < EQL | | | | | ug/L |
| | Mercury, total recoverable | <.37 | U | V | < EQL | 1.47 | | | < 2 | ug/L |
| | Methacrylonitrile | <200 | U | | < EQL | <200 | U | | < EQL | ug/L |
| | Methapyrilene | <11 | JU | Q | < EQL | | | | | ug/L |
| | Methoxychlor | <1 | U | | < EQL | | | | | ug/L |
| | Methyl ethyl ketone | <20 | U | | < EQL | <20 | U | | < EQL | ug/L |
| | Methyl isobutyl ketone | <10 | U | | < EQL | <10 | U | | < EQL | ug/L |
| | Methyl methacrylate | <20 | U | | < EQL | <20 | U | | < EQL | ug/L |
| | Methyl methanesulfonate | <11 | JU | Q | < EQL | | | | | ug/L |
| | N-Nitrosodi-n-butylamine | <11 | JU | Q | < EQL | | | | | ug/L |
| | N-Nitrosodiethylamine | <11 | JU | Q | < EQL | | | | | ug/L |
| | N-Nitrosodimethylamine | <26 | JU | Q | < EQL | | | | | ug/L |
| | N-Nitrosodiphenylamine | <11 | JU | Q | < EQL | | | | | ug/L |
| | N-Nitrosodipropylamine | <11 | JU | Q | < EQL | | | | | ug/L |
| | N-Nitrosomethylethylamine | <11 | JU | Q | < EQL | | | | | ug/L |
| | N-Nitrosopiperidine | <11 | JU | Q | < EQL | | | | | ug/L |
| | N-Nitrosopyrrolidine | <11 | JU | Q | < EQL | | | | | ug/L |
| | Naphthalene | <11 | JU | Q | < EQL | | | | | ug/L |
| | Nickel, total recoverable | <50 | U | | < EQL | <50 | U | | < EQL | ug/L |
| | Nitrobenzene | <11 | JU | Q | < EQL | | | | | ug/L |
| | O,O,O-Triethyl phosphorothioate | <11 | JU | Q | < EQL | | | | | ug/L |
| | PCB | | | | | | | | | |
| | PCB 1016 | <1 | U | | < EQL | | | | | ug/L |
| | PCB 1232 | <1 | U | | < EQL | | | | | ug/L |
| | PCB 1242 | <2.1 | U | | < EQL | | | | | ug/L |
| | PCB 1248 | <1 | U | | < EQL | | | | | ug/L |
| | PCB 1254 | <1 | U | | < EQL | | | | | ug/L |
| | PCB 1260 | <1 | U | | < EQL | | | | | ug/L |
| | Parathion | <11 | JU | Q | < EQL | | | | | ug/L |
| | Parathion methyl | <11 | JU | Q | < EQL | | | | | ug/L |
| | Pentachlorobenzene | <53 | JU | Q | < EQL | | | | | ug/L |
| | Pentachloronitrobenzene | <11 | JU | Q | < EQL | | | | | ug/L |
| | Pentachlorophenol | <26 | JU | Q | < EQL | | | | | ug/L |
| | Phenacetin | <11 | JU | Q | < EQL | | | | | ug/L |
| | Phenanthrene | <11 | JU | Q | < EQL | | | | | ug/L |
| | Phenol | <11 | JU | Q | < EQL | | | | | ug/L |
| | Phorate | <11 | JU | Q | < EQL | | | | | ug/L |
| | Pronamid | <11 | JU | Q | < EQL | | | | | ug/L |
| | Propionitrile | <200 | U | | < EQL | <200 | U | | < EQL | ug/L |
| | Pyrene | <11 | JU | Q | < EQL | | | | | ug/L |
| | Safrole | <11 | JU | Q | < EQL | | | | | ug/L |
| | Selenium, total recoverable | <10 | U | | < EQL | <10 | U | | < EQL | ug/L |
| | Silver, total recoverable | <20 | U | | < EQL | <5 | JU | | < EQL | ug/L |
| | Specific conductance | 25 | | | < 37 | 28.5 | | | < 37 | uS/cm |
| | Styrene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Sulfide | 1000 | R | Q | Rejected | | | | | ug/L |

Notes: Concentrations in bold italics exceed the groundwater protection standards listed in Appendix A. Bold italics were not assigned if the result was qualified with a "J", "R", "L", or "U" qualifier.

+ = exceeded the Groundwater Protection Standards listed in Appendix A.

Table B. Assessment Monitoring Results for Groundwater Wells (Cont.)

WELL LFW 32

ANALYTICAL DATA

Groundwater Protection Standard

| ST | Parameter | 1st Half | CLP | EPA | Filt. | 2nd Half | CLP | EPA | Filt. | Unit |
|-----------|-----------------------------|-----------------|------------|------------|--------------|-----------------|------------|------------|--------------|-------------|
| | Tetrachloroethylene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Thallium, total recoverable | <10 | U | | < EQL | 7.77 | J | I | NDD | ug/L |
| | Thionazin | <11 | JU | Q | < EQL | | | | | ug/L |
| | Tin, total recoverable | <200 | U | | < EQL | | | | | ug/L |
| | Toluene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Toxaphene | <2.1 | U | | < EQL | | | | | ug/L |
| | Trichloroethylene | .91 | J | I | NDD | 1.2 | J | IK | NDD | ug/L |
| | Trichlorofluoromethane | 160 | | | > 20 | 200 | J | K | NDD | ug/L |
| | Vanadium, total recoverable | <10 | U | | < EQL | <10 | U | | < EQL | ug/L |
| | Vinyl acetate | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Xylenes | <10 | U | | < EQL | <10 | U | | < EQL | ug/L |
| | Zinc, total recoverable | <7.99 | JU | | < EQL | <20 | U | | < EQL | ug/L |
| | alpha-Benzene hexachloride | <.1 | U | | < EQL | | | | | ug/L |
| | alpha-Chlordane | <.1 | U | | < EQL | | | | | ug/L |
| | beta-Benzene hexachloride | <.1 | U | | < EQL | <.095 | U | | < EQL | ug/L |
| | cis-1,2-Dichloroethylene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | cis-1,3-Dichloropropene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | delta-Benzene hexachloride | <.1 | U | | < EQL | | | | | ug/L |
| | gamma-Chlordane | <.1 | U | | < EQL | | | | | ug/L |
| | m-Cresol (3-Methylphenol) | | | | | | | | | |
| | m-Nitroaniline | <26 | JU | Q | < EQL | | | | | ug/L |
| | o-Cresol (2-Methylphenol) | <11 | JU | Q | < EQL | | | | | ug/L |
| | o-Nitroaniline | <26 | JU | Q | < EQL | | | | | ug/L |
| | o-Toluidine | <11 | JU | Q | < EQL | | | | | ug/L |
| | p,p"-DDD | <.21 | U | | < EQL | | | | | ug/L |
| | p,p"-DDE | <.21 | U | | < EQL | | | | | ug/L |
| | p,p"-DDT | <.21 | U | | < EQL | | | | | ug/L |
| | p-Cresol (4-Methylphenol) | <10 | JU | Q | < EQL | | | | | ug/L |
| | p-Dimethylaminoazobenzene | <11 | JU | Q | < EQL | | | | | ug/L |
| | p-Nitroaniline | <11 | JU | Q | < EQL | | | | | ug/L |
| | p-Phenylenediamine | <11 | JU | Q | < EQL | | | | | ug/L |
| | pH | 5.8 | J | Q | NDD | 4.88 | J | Q | NDD | pH |
| | trans-1,2-Dichloroethylene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | trans-1,3-Dichloropropene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | trans-1,4-Dichloro-2-butene | <20 | U | | < EQL | <20 | U | | < EQL | ug/L |

Notes: Concentrations in bold italics exceed the groundwater protection standards listed in Appendix A. Bold italics were not assigned if the result was qualified with a "J", "R", "L", or "U" qualifier.

+ = exceeded the Groundwater Protection Standards listed in Appendix A.

Table B. Assessment Monitoring Results for Groundwater Wells (Cont.)

WELL LFW 32C

| <u>SRS Coord.</u> | <u>Lat/Longitude</u> | <u>Screen Zone Elevation</u> | <u>Top of Casing</u> | <u>Casing</u> | <u>Pump</u> | <u>Screen Zone</u> |
|----------------------|----------------------------------|------------------------------|----------------------|---------------|-------------|--------------------|
| N 85837.8 E 44923 | 33.28891 Deg N -81.7152 Deg W | 113.6 - 98.6 ft msl | 223.6 ft msl | 2" PVC | V | Middle |

SAMPLE DATE 02/28/00 10/02/00

FIELD DATA

| <u>Parameter</u> | <u>1st Half</u> | <u>2nd Half</u> | <u>Unit</u> |
|---------------------|-----------------|-----------------|-------------|
| Water Elevation | 158.61 | 157.89 | ft msl |
| pH | 5 | 5.4 | pH |
| Sp. Conductance | 28 | 29 | uS/cm |
| Water temperature | 18.1 | 19.2 | deg. C |
| Alkalinity as CaCO3 | 1 | 0 | mg/L |
| Turbidity | .9 | .5 | NTU |
| Volumes purged | 2.98 | 4.73 | well volume |
| Sampling code | | | |

ANALYTICAL DATA

Groundwater Protection Standard

| <u>SI</u> | <u>Parameter</u> | <u>1st Half</u> | <u>CLP</u> | <u>EPA</u> | <u>Filt.</u> | <u>2nd Half</u> | <u>CLP</u> | <u>EPA</u> | <u>Filt.</u> | <u>Unit</u> |
|-----------|--------------------------------|-----------------|------------|------------|--------------|-----------------|------------|------------|--------------|-------------|
| | 1,1,1,2-Tetrachloroethane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,1,1-Trichloroethane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,1,2,2-Tetrachloroethane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,1,2-Trichloroethane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,1-Dichloroethane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,1-Dichloroethylene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,1-Dichloropropene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,2,3-Trichloropropane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,2,4,5-Tetrachlorobenzene | <52 | JU | L | < EQL | | | | | ug/L |
| | 1,2,4-Trichlorobenzene | <11 | JU | L | < EQL | | | | | ug/L |
| | 1,2-Dibromo-3-chloropropane | <10 | U | | < EQL | <10 | U | | < EQL | ug/L |
| | 1,2-Dibromoethane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,2-Dichlorobenzene | <11 | JU | L | < EQL | <5 | U | | < EQL | ug/L |
| | 1,2-Dichloroethane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,2-Dichloropropane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,3,5-Trinitrobenzene | <10 | JU | L | < EQL | | | | | ug/L |
| | 1,3-Dichlorobenzene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,3-Dichloropropane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,3-Dinitrobenzene | <50 | JU | L | < EQL | | | | | ug/L |
| | 1,4-Dichlorobenzene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,4-Naphthoquinone | <11 | JU | L | < EQL | | | | | ug/L |
| | 1-Naphthylamine | <11 | JU | L | < EQL | | | | | ug/L |
| | 2,2-Dichloropropane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 2,2-Oxybis(1-chloropropane) | <11 | JU | L | < EQL | | | | | ug/L |
| | 2,3,4,6-Tetrachlorophenol | <11 | JU | L | < EQL | | | | | ug/L |
| | 2,4,5-T | <2 | U | | < EQL | | | | | ug/L |
| | 2,4,5-TP (Silvex) | <2 | U | | < EQL | | | | | ug/L |
| | 2,4,5-Trichlorophenol | <11 | JU | L | < EQL | | | | | ug/L |
| | 2,4,6-Trichlorophenol | <26 | JU | L | < EQL | | | | | ug/L |
| | 2,4-Dichlorophenol | <11 | JU | L | < EQL | | | | | ug/L |
| | 2,4-Dichlorophenoxyacetic acid | <2 | U | | < EQL | | | | | ug/L |
| | 2,4-Dimethyl phenol | <11 | JU | L | < EQL | | | | | ug/L |
| | 2,4-Dinitrophenol | <26 | JU | L | < EQL | | | | | ug/L |
| | 2,4-Dinitrotoluene | <11 | JU | L | < EQL | | | | | ug/L |
| | 2,6-Dichlorophenol | <25 | JU | L | < EQL | | | | | ug/L |
| | 2,6-Dinitrotoluene | <11 | JU | L | < EQL | | | | | ug/L |
| | 2-Acetylaminofluorene | <11 | JU | L | < EQL | | | | | ug/L |
| | 2-Chloronaphthalene | <11 | JU | L | < EQL | | | | | ug/L |
| | 2-Chlorophenol | <11 | JU | L | < EQL | | | | | ug/L |
| | 2-Hexanone | <20 | U | | < EQL | <20 | U | | < EQL | ug/L |
| | 2-Methyl-4,6-dinitrophenol | <26 | JU | L | < EQL | | | | | ug/L |
| | 2-Methylnaphthalene | <11 | JU | L | < EQL | | | | | ug/L |
| | 2-Naphthylamine | <11 | JU | L | < EQL | | | | | ug/L |
| | 2-Nitrophenol | <11 | JU | L | < EQL | | | | | ug/L |
| | 2-sec-Butyl-4,6-dinitrophenol | <2 | JU | L | < EQL | | | | | ug/L |
| | 3,3"-Dichlorobenzidine | <11 | JU | L | < EQL | | | | | ug/L |
| | 3,3"-Dimethylbenzidine | <21 | JU | L | < EQL | | | | | ug/L |
| | 3-Methylcholanthrene | <11 | JU | L | < EQL | | | | | ug/L |
| | 4-Aminobiphenyl | <11 | JU | L | < EQL | | | | | ug/L |
| | 4-Bromophenyl phenyl ether | <11 | JU | L | < EQL | | | | | ug/L |
| | 4-Chloro-m-cresol | <11 | JU | L | < EQL | | | | | ug/L |

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+ = exceeded the Groundwater Protection Standards listed in Appendix A.

Table B. Assessment Monitoring Results for Groundwater Wells (Cont.)

WELL LFW 32C

ANALYTICAL DATA

Groundwater Protection Standard

| ST | Parameter | 1st Half | CLP | EPA | Filt. | 2nd Half | CLP | EPA | Filt. | Unit |
|----|--------------------------------------|----------|-----|-----|-------|----------|-----|-----|-------|------|
| | 4-Chloroaniline | <11 | JU | L | < EQL | | | | | ug/L |
| | 4-Chlorophenyl phenyl ether | <11 | JU | L | < EQL | | | | | ug/L |
| | 4-Nitrophenol | <26 | JU | L | < EQL | | | | | ug/L |
| | 5-Nitro-o-toluidine | <11 | JU | L | < EQL | | | | | ug/L |
| | 7,12-Dimethylbenz(a)anthracene | <11 | JU | L | < EQL | | | | | ug/L |
| | Acenaphthene | <11 | JU | L | < EQL | | | | | ug/L |
| | Acenaphthylene | <11 | JU | L | < EQL | | | | | ug/L |
| | Acetone | <20 | U | | < EQL | <20 | U | | < EQL | ug/L |
| | Acetonitrile (Methyl cyanide) | <200 | U | | < EQL | <200 | U | | < EQL | ug/L |
| | Acetophenone | <11 | JU | L | < EQL | | | | | ug/L |
| | Acrolein | <50 | U | | < EQL | <50 | U | | < EQL | ug/L |
| | Acrylonitrile | <10 | U | | < EQL | <10 | U | | < EQL | ug/L |
| | Aldrin | <.11 | U | | < EQL | | | | | ug/L |
| | Allyl chloride | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Anthracene | <11 | JU | L | < EQL | | | | | ug/L |
| | Antimony, total recoverable | <100 | U | | < EQL | <100 | U | | < EQL | ug/L |
| | Arsenic, total recoverable | <10 | U | | < EQL | <10 | U | | < EQL | ug/L |
| | Barium, total recoverable | 5.78 | J | I | NDD | 5.83 | J | I | NDD | ug/L |
| | Benzene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Benzo(a)anthracene | <11 | JU | L | < EQL | | | | | ug/L |
| | Benzo(a)pyrene | <11 | JU | L | < EQL | | | | | ug/L |
| | Benzo(b)fluoranthene | <11 | JU | L | < EQL | | | | | ug/L |
| | Benzo(g,h,i)perylene | <11 | JU | L | < EQL | | | | | ug/L |
| | Benzo(k)fluoranthene | <11 | JU | L | < EQL | | | | | ug/L |
| | Benzyl alcohol | | | | | | | | | |
| | Beryllium, total recoverable | <.14 | JU | V | < EQL | <.19 | JU | | < EQL | ug/L |
| | Bis(2-chloroethoxy) methane | <11 | JU | L | < EQL | | | | | ug/L |
| | Bis(2-chloroethyl) ether | <11 | JU | L | < EQL | | | | | ug/L |
| | Bis(2-ethylhexyl) phthalate | <11 | JU | L | < EQL | | | | | ug/L |
| | Bromochloromethane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Bromodichloromethane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Bromoform | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Bromomethane (Methyl bromide) | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Butylbenzyl phthalate | <11 | JU | L | < EQL | | | | | ug/L |
| | Cadmium, total recoverable | <10 | U | | < EQL | <10 | U | | < EQL | ug/L |
| | Carbon disulfide | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Carbon tetrachloride | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Chlorobenzene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Chlorobenzilate | <11 | JU | L | < EQL | | | | | ug/L |
| | Chloroethane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Chloroethene (Vinyl chloride) | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Chloroform | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Chloromethane (Methyl chloride) | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Chloroprene | <20 | U | | < EQL | <20 | U | | < EQL | ug/L |
| | Chromium, total recoverable | <10 | U | | < EQL | <10 | U | | < EQL | ug/L |
| | Chrysene | <11 | JU | L | < EQL | | | | | ug/L |
| | Cobalt, total recoverable | <20 | U | | < EQL | <20 | U | | < EQL | ug/L |
| | Copper, total recoverable | <20 | U | | < EQL | <20 | U | | < EQL | ug/L |
| | Cyanide | <10 | U | | < EQL | | | | | ug/L |
| | Di-n-butyl phthalate | <11 | JU | L | < EQL | <9.6 | U | | < EQL | ug/L |
| | Di-n-octyl phthalate | <11 | JU | L | < EQL | | | | | ug/L |
| | Diallate | <11 | JU | L | < EQL | | | | | ug/L |
| | Dibenz(a,h)anthracene | <11 | JU | L | < EQL | | | | | ug/L |
| | Dibenzofuran | <11 | JU | L | < EQL | | | | | ug/L |
| | Dibromochloromethane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Dibromomethane (Methylene bromide) | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Dichlorodifluoromethane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Dichloromethane (Methylene chloride) | <2.7 | U | V | < EQL | <10 | U | | < EQL | ug/L |
| | Dieldrin | <.21 | U | | < EQL | | | | | ug/L |
| | Diethyl phthalate | <11 | JU | L | < EQL | | | | | ug/L |
| | Dimethoate | <11 | JU | L | < EQL | | | | | ug/L |
| | Dimethyl phthalate | <11 | JU | L | < EQL | | | | | ug/L |
| | Diphenylamine | <11 | JU | L | < EQL | | | | | ug/L |
| | Disulfoton | <11 | JU | L | < EQL | | | | | ug/L |
| | Endosulfan I | <.11 | U | | < EQL | | | | | ug/L |
| | Endosulfan II | <.21 | U | | < EQL | | | | | ug/L |
| | Endosulfan sulfate | <.21 | U | | < EQL | | | | | ug/L |
| | Endrin | <.21 | U | | < EQL | | | | | ug/L |

Notes: Concentrations in bold italics exceed the groundwater protection standards listed in Appendix A. Bold italics were not assigned if the result was qualified with a "J", "R", "L", or "U" qualifier.

+ = exceeded the Groundwater Protection Standards listed in Appendix A.

Table B. Assessment Monitoring Results for Groundwater Wells (Cont.)

WELL LFW 32C

ANALYTICAL DATA

Groundwater Protection Standard

| ST | Parameter | 1st Half | CLP | EPA | Filt. | 2nd Half | CLP | EPA | Filt. | Unit |
|----|---------------------------------|----------|-----|-----|-------|----------|-----|-----|-------|-------|
| | Endrin aldehyde | <.21 | U | | < EQL | | | | | ug/L |
| | Ethyl methacrylate | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Ethyl methanesulfonate | <11 | JU | L | < EQL | | | | | ug/L |
| | Ethylbenzene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Famphur | <200 | JU | L | < EQL | | | | | ug/L |
| | Fluoranthene | <11 | JU | L | < EQL | | | | | ug/L |
| | Fluorene | <11 | JU | L | < EQL | | | | | ug/L |
| | Heptachlor | <.11 | U | | < EQL | | | | | ug/L |
| | Heptachlor epoxide | <.11 | U | | < EQL | | | | | ug/L |
| | Hexachlorobenzene | <11 | JU | L | < EQL | | | | | ug/L |
| | Hexachlorobutadiene | <11 | JU | L | < EQL | | | | | ug/L |
| | Hexachlorocyclopentadiene | <11 | JU | L | < EQL | | | | | ug/L |
| | Hexachloroethane | <11 | JU | L | < EQL | | | | | ug/L |
| | Hexachloropropene | <52 | JU | L | < EQL | | | | | ug/L |
| | Indeno(1,2,3-c,d)pyrene | <11 | JU | L | < EQL | | | | | ug/L |
| | Iodomethane (Methyl iodide) | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Isobutyl alcohol | <500 | U | | < EQL | <500 | U | | < EQL | ug/L |
| | Isodrin | <11 | JU | L | < EQL | | | | | ug/L |
| | Isophorone | <11 | JU | L | < EQL | | | | | ug/L |
| | Isosafrole | <11 | JU | L | < EQL | | | | | ug/L |
| | Kepone | <11 | JU | L | < EQL | | | | | ug/L |
| | Lead, total recoverable | <10 | U | | < EQL | <10 | U | | < EQL | ug/L |
| | Lindane | <.11 | U | | < EQL | | | | | ug/L |
| | Mercury, total recoverable | <.5 | U | | < EQL | <.5 | U | | < EQL | ug/L |
| | Methacrylonitrile | <200 | U | | < EQL | <200 | U | | < EQL | ug/L |
| | Methapyrilene | <11 | JU | L | < EQL | | | | | ug/L |
| | Methoxychlor | <1.1 | U | | < EQL | | | | | ug/L |
| | Methyl ethyl ketone | <20 | U | | < EQL | <20 | U | | < EQL | ug/L |
| | Methyl isobutyl ketone | <10 | U | | < EQL | <10 | U | | < EQL | ug/L |
| | Methyl methacrylate | <20 | U | | < EQL | <20 | U | | < EQL | ug/L |
| | Methyl methanesulfonate | <11 | JU | L | < EQL | | | | | ug/L |
| | N-Nitrosodi-n-butylamine | <11 | JU | L | < EQL | | | | | ug/L |
| | N-Nitrosodiethylamine | <11 | JU | L | < EQL | | | | | ug/L |
| | N-Nitrosodimethylamine | <26 | JU | L | < EQL | | | | | ug/L |
| | N-Nitrosodiphenylamine | <11 | JU | L | < EQL | | | | | ug/L |
| | N-Nitrosodipropylamine | <11 | JU | L | < EQL | | | | | ug/L |
| | N-Nitrosomethylethylamine | <11 | JU | L | < EQL | | | | | ug/L |
| | N-Nitrosopiperidine | <11 | JU | L | < EQL | | | | | ug/L |
| | N-Nitrosopyrrolidine | <11 | JU | L | < EQL | | | | | ug/L |
| | Naphthalene | <11 | JU | L | < EQL | | | | | ug/L |
| | Nickel, total recoverable | <50 | U | | < EQL | <50 | U | | < EQL | ug/L |
| | Nitrobenzene | <11 | JU | L | < EQL | | | | | ug/L |
| | O,O,O-Triethyl phosphorothioate | <11 | JU | L | < EQL | | | | | ug/L |
| | PCB | | | | | | | | | |
| | PCB 1016 | <1.1 | U | | < EQL | | | | | ug/L |
| | PCB 1232 | <1.1 | U | | < EQL | | | | | ug/L |
| | PCB 1242 | <2.1 | U | | < EQL | | | | | ug/L |
| | PCB 1248 | <1.1 | U | | < EQL | | | | | ug/L |
| | PCB 1254 | <1.1 | U | | < EQL | | | | | ug/L |
| | PCB 1260 | <1.1 | U | | < EQL | | | | | ug/L |
| | Parathion | <11 | JU | L | < EQL | | | | | ug/L |
| | Parathion methyl | <11 | JU | L | < EQL | | | | | ug/L |
| | Pentachlorobenzene | <52 | JU | L | < EQL | | | | | ug/L |
| | Pentachloronitrobenzene | <11 | JU | L | < EQL | | | | | ug/L |
| | Pentachlorophenol | <26 | JU | L | < EQL | | | | | ug/L |
| | Phenacetin | <11 | JU | L | < EQL | | | | | ug/L |
| | Phenanthrene | <11 | JU | L | < EQL | | | | | ug/L |
| | Phenol | <11 | JU | L | < EQL | | | | | ug/L |
| | Phorate | <11 | JU | L | < EQL | | | | | ug/L |
| | Pronamid | <11 | JU | L | < EQL | | | | | ug/L |
| | Propionitrile | <200 | U | | < EQL | <200 | U | | < EQL | ug/L |
| | Pyrene | <11 | JU | L | < EQL | | | | | ug/L |
| | Safrole | <11 | JU | L | < EQL | | | | | ug/L |
| | Selenium, total recoverable | <10 | U | | < EQL | <10 | U | | < EQL | ug/L |
| | Silver, total recoverable | <20 | U | | < EQL | <20 | U | | < EQL | ug/L |
| | Specific conductance | 27.7 | | | < 37 | 27.7 | | | < 37 | uS/cm |
| | Styrene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Sulfide | <1000 | U | | < EQL | | | | | ug/L |

Notes: Concentrations in bold italics exceed the groundwater protection standards listed in Appendix A. Bold italics were not assigned if the result was qualified with a "J", "R", "L", or "U" qualifier.

+ = exceeded the Groundwater Protection Standards listed in Appendix A.

Table B. Assessment Monitoring Results for Groundwater Wells (Cont.)

WELL LFW 32C

ANALYTICAL DATA

Groundwater Protection Standard

| ST | Parameter | 1st Half | CLP | EPA | Filt. | 2nd Half | CLP | EPA | Filt. | Unit |
|----|-----------------------------|----------|-----|-----|-------|----------|-----|-----|-------|------|
| | Tetrachloroethylene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Thallium, total recoverable | <10 | U | | < EQL | <5.39 | JU | | < EQL | ug/L |
| | Thionazin | <11 | JU | L | < EQL | | | | | ug/L |
| | Tin, total recoverable | <200 | U | | < EQL | | | | | ug/L |
| | Toluene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Toxaphene | <2.1 | U | | < EQL | | | | | ug/L |
| | Trichloroethylene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Trichlorofluoromethane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Vanadium, total recoverable | <10 | U | | < EQL | <10 | U | | < EQL | ug/L |
| | Vinyl acetate | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Xylenes | <10 | U | | < EQL | <10 | U | | < EQL | ug/L |
| | Zinc, total recoverable | <20 | U | | < EQL | <20 | U | | < EQL | ug/L |
| | alpha-Benzene hexachloride | <.11 | U | | < EQL | | | | | ug/L |
| | alpha-Chlordane | <.11 | U | | < EQL | | | | | ug/L |
| | beta-Benzene hexachloride | <.11 | U | | < EQL | <.095 | U | | < EQL | ug/L |
| | cis-1,2-Dichloroethylene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | cis-1,3-Dichloropropene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | delta-Benzene hexachloride | <.11 | U | | < EQL | | | | | ug/L |
| | gamma-Chlordane | <.11 | U | | < EQL | | | | | ug/L |
| | m-Cresol (3-Methylphenol) | | | | | | | | | |
| | m-Nitroaniline | <26 | JU | L | < EQL | | | | | ug/L |
| | o-Cresol (2-Methylphenol) | <11 | JU | L | < EQL | | | | | ug/L |
| | o-Nitroaniline | <26 | JU | L | < EQL | | | | | ug/L |
| | o-Toluidine | <11 | JU | L | < EQL | | | | | ug/L |
| | p,p"-DDD | <.21 | U | | < EQL | | | | | ug/L |
| | p,p"-DDE | <.21 | U | | < EQL | | | | | ug/L |
| | p,p"-DDT | <.21 | U | | < EQL | | | | | ug/L |
| | p-Cresol (4-Methylphenol) | <10 | JU | L | < EQL | | | | | ug/L |
| | p-Dimethylaminoazobenzene | <11 | JU | L | < EQL | | | | | ug/L |
| | p-Nitroaniline | <11 | JU | L | < EQL | | | | | ug/L |
| | p-Phenylenediamine | <11 | JU | L | < EQL | | | | | ug/L |
| | pH | 5.55 | J | Q | NDD | 4.92 | J | Q | NDD | pH |
| | trans-1,2-Dichloroethylene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | trans-1,3-Dichloropropene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | trans-1,4-Dichloro-2-butene | <20 | U | | < EQL | <20 | U | | < EQL | ug/L |

Notes: Concentrations in bold italics exceed the groundwater protection standards listed in Appendix A. Bold italics were not assigned if the result was qualified with a "J", "R", "L", or "U" qualifier.

+ = exceeded the Groundwater Protection Standards listed in Appendix A.

Table B. Assessment Monitoring Results for Groundwater Wells (Cont.)

WELL LFW 34

| <u>SRS Coord.</u> | <u>Lat/Longitude</u> | <u>Screen Zone Elevation</u> | <u>Top of Casing</u> | <u>Casing</u> | <u>Pump</u> | <u>Screen Zone</u> |
|------------------------|----------------------------------|------------------------------|----------------------|---------------|-------------|--------------------|
| N 85409.5 E 45016.9 | 33.28811 Deg N -81.7142 Deg W | 164.7 - 143.7 ft msl | 201 ft msl | 4 " PVC | S | Upper |

SAMPLE DATE 03/08/00 08/17/00

FIELD DATA

| <u>Parameter</u> | <u>1st Half</u> | <u>2nd Half</u> | <u>Unit</u> |
|---------------------|-----------------|-----------------|-------------|
| Water Elevation | 158.6 | 146.65 | ft msl |
| pH | 5 | 5.1 | pH |
| Sp. Conductance | 31 | 38 | uS/cm |
| Water temperature | 19.5 | 23.2 | deg. C |
| Alkalinity as CaCO3 | 0 | 0 | mg/L |
| Turbidity | .2 | .6 | NTU |
| Volumes purged | 5.60 | 7.76 | well volume |
| Sampling code | | | |

ANALYTICAL DATA

Groundwater Protection Standard

| <u>ST</u> | <u>Parameter</u> | <u>1st Half</u> | <u>CLP</u> | <u>EPA</u> | <u>Filt.</u> | <u>2nd Half</u> | <u>CLP</u> | <u>EPA</u> | <u>Filt.</u> | <u>Unit</u> |
|-----------|--------------------------------|-----------------|------------|------------|--------------|-----------------|------------|------------|--------------|-------------|
| | 1,1,1,2-Tetrachloroethane | <25 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,1,1-Trichloroethane | 14 | | | < 200 | 15 | | | < 200 | ug/L |
| | 1,1,2,2-Tetrachloroethane | <25 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,1,2-Trichloroethane | <25 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,1-Dichloroethane | 8.7 | | | < 20 | 14 | | | < 20 | ug/L |
| | 1,1-Dichloroethylene | <25 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,1-Dichloropropene | <25 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,2,3-Trichloropropane | <25 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,2,4,5-Tetrachlorobenzene | <48 | JU | Q | < EQL | | | | | ug/L |
| | 1,2,4-Trichlorobenzene | <9.6 | JU | Q | < EQL | | | | | ug/L |
| | 1,2-Dibromo-3-chloropropane | <50 | U | | < EQL | <10 | U | | < EQL | ug/L |
| | 1,2-Dibromoethane | <25 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,2-Dichlorobenzene | <25 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,2-Dichloroethane | <25 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,2-Dichloropropane | <25 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,3,5-Trinitrobenzene | <10 | JU | Q | < EQL | | | | | ug/L |
| | 1,3-Dichlorobenzene | <25 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,3-Dichloropropane | <25 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,3-Dinitrobenzene | <50 | JU | Q | < EQL | | | | | ug/L |
| | 1,4-Dichlorobenzene | <25 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,4-Naphthoquinone | <9.6 | JU | Q | < EQL | | | | | ug/L |
| | 1-Naphthylamine | <9.6 | JU | Q | < EQL | | | | | ug/L |
| | 2,2-Dichloropropane | <25 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 2,2-Oxybis(1-chloropropane) | <9.6 | JU | Q | < EQL | | | | | ug/L |
| | 2,3,4,6-Tetrachlorophenol | <9.6 | JU | Q | < EQL | | | | | ug/L |
| | 2,4,5-T | <.2 | U | Q | < EQL | | | | | ug/L |
| | 2,4,5-TP (Silvex) | <.2 | U | Q | < EQL | | | | | ug/L |
| | 2,4,5-Trichlorophenol | <9.6 | JU | Q | < EQL | | | | | ug/L |
| | 2,4,6-Trichlorophenol | <24 | JU | Q | < EQL | | | | | ug/L |
| | 2,4-Dichlorophenol | <9.6 | JU | Q | < EQL | | | | | ug/L |
| | 2,4-Dichlorophenoxyacetic acid | <.2 | U | Q | < EQL | | | | | ug/L |
| | 2,4-Dimethyl phenol | <9.6 | JU | Q | < EQL | | | | | ug/L |
| | 2,4-Dinitrophenol | <24 | JU | Q | < EQL | | | | | ug/L |
| | 2,4-Dinitrotoluene | <9.6 | JU | Q | < EQL | | | | | ug/L |
| | 2,6-Dichlorophenol | <25 | JU | Q | < EQL | | | | | ug/L |
| | 2,6-Dinitrotoluene | <9.6 | JU | Q | < EQL | | | | | ug/L |
| | 2-Acetylaminofluorene | <9.6 | JU | Q | < EQL | | | | | ug/L |
| | 2-Chloronaphthalene | <9.6 | JU | Q | < EQL | | | | | ug/L |
| | 2-Chlorophenol | <9.6 | JU | Q | < EQL | | | | | ug/L |
| | 2-Hexanone | <100 | U | | < EQL | <20 | U | | < EQL | ug/L |
| | 2-Methyl-4,6-dinitrophenol | <24 | JU | Q | < EQL | | | | | ug/L |
| | 2-Methylnaphthalene | <9.6 | JU | Q | < EQL | | | | | ug/L |
| | 2-Naphthylamine | <9.6 | JU | Q | < EQL | | | | | ug/L |
| | 2-Nitrophenol | <9.6 | JU | Q | < EQL | | | | | ug/L |
| | 2-sec-Butyl-4,6-dinitrophenol | <.2 | U | Q | < EQL | | | | | ug/L |
| | 3,3"-Dichlorobenzidine | <9.6 | JU | Q | < EQL | | | | | ug/L |
| | 3,3"-Dimethylbenzidine | <19 | JU | Q | < EQL | | | | | ug/L |
| | 3-Methylcholanthrene | <9.6 | JU | Q | < EQL | | | | | ug/L |
| | 4-Aminobiphenyl | <9.6 | JU | Q | < EQL | | | | | ug/L |
| | 4-Bromophenyl phenyl ether | <9.6 | JU | Q | < EQL | | | | | ug/L |
| | 4-Chloro-m-cresol | <9.6 | JU | Q | < EQL | | | | | ug/L |

Notes: Concentrations in bold italics exceed the groundwater protection standards listed in Appendix A. Bold italics were not assigned if the result was qualified with a "J", "R", "L", or "U" qualifier.

+ = exceeded the Groundwater Protection Standards listed in Appendix A.

Table B. Assessment Monitoring Results for Groundwater Wells (Cont.)

WELL LFW 34

ANALYTICAL DATA

Groundwater Protection Standard

| ST | Parameter | 1st Half | CLP | EPA | Filt. | 2nd Half | CLP | EPA | Filt. | Unit |
|----|--------------------------------------|----------|-----|-----|----------|----------|-----|-----|--------|------|
| | 4-Chloroaniline | <9.6 | JU | Q | < EQL | | | | | ug/L |
| | 4-Chlorophenyl phenyl ether | <9.6 | JU | Q | < EQL | | | | | ug/L |
| | 4-Nitrophenol | <24 | JU | Q | < EQL | | | | | ug/L |
| | 5-Nitro-o-toluidine | <9.6 | JU | Q | < EQL | | | | | ug/L |
| | 7,12-Dimethylbenz(a)anthracene | <9.6 | JU | Q | < EQL | | | | | ug/L |
| | Acenaphthene | <9.6 | JU | Q | < EQL | | | | | ug/L |
| | Acenaphthylene | <9.6 | JU | Q | < EQL | | | | | ug/L |
| | Acetone | <100 | U | | < EQL | <20 | U | | < EQL | ug/L |
| | Acetonitrile (Methyl cyanide) | <1000 | U | | < EQL | <200 | U | | < EQL | ug/L |
| | Acetophenone | <9.6 | JU | Q | < EQL | | | | | ug/L |
| | Acrolein | <250 | U | | < EQL | <50 | U | | < EQL | ug/L |
| | Acrylonitrile | <50 | U | | < EQL | <10 | U | | < EQL | ug/L |
| | Aldrin | <.1 | U | | < EQL | | | | | ug/L |
| | Allyl chloride | <25 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Anthracene | <9.6 | JU | Q | < EQL | | | | | ug/L |
| | Antimony, total recoverable | <100 | U | | < EQL | <100 | U | | < EQL | ug/L |
| | Arsenic, total recoverable | <10 | U | | < EQL | <10 | U | | < EQL | ug/L |
| | Barium, total recoverable | 16.8 | | | < 2000 | 14.5 | | | < 2000 | ug/L |
| | Benzene | <25 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Benzo(a)anthracene | <9.6 | JU | Q | < EQL | | | | | ug/L |
| | Benzo(a)pyrene | <9.6 | JU | Q | < EQL | | | | | ug/L |
| | Benzo(b)fluoranthene | <9.6 | JU | Q | < EQL | | | | | ug/L |
| | Benzo(g,h,i)perylene | <9.6 | JU | Q | < EQL | | | | | ug/L |
| | Benzo(k)fluoranthene | <9.6 | JU | Q | < EQL | | | | | ug/L |
| | Benzyl alcohol | | | | | | | | | |
| | Beryllium, total recoverable | <1 | U | | < EQL | <1 | U | | < EQL | ug/L |
| | Bis(2-chloroethoxy) methane | <9.6 | JU | Q | < EQL | | | | | ug/L |
| | Bis(2-chloroethyl) ether | <9.6 | JU | Q | < EQL | | | | | ug/L |
| | Bis(2-ethylhexyl) phthalate | <9.6 | JU | Q | < EQL | | | | | ug/L |
| | Bromochloromethane | <25 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Bromodichloromethane | <25 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Bromoform | <25 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Bromomethane (Methyl bromide) | <25 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Butylbenzyl phthalate | <9.6 | JU | Q | < EQL | | | | | ug/L |
| | Cadmium, total recoverable | <10 | U | | < EQL | <10 | U | | < EQL | ug/L |
| | Carbon disulfide | <25 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Carbon tetrachloride | <25 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Chlorobenzene | <25 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Chlorobenzilate | <9.6 | JU | Q | < EQL | | | | | ug/L |
| | Chloroethane | <25 | U | | < EQL | 4.2 | J | I | NDD | ug/L |
| | Chloroethene (Vinyl chloride) | <25 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Chloroform | <25 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Chloromethane (Methyl chloride) | <25 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Chloroprene | <100 | U | | < EQL | <20 | U | | < EQL | ug/L |
| | Chromium, total recoverable | <10 | U | | < EQL | <10 | U | | < EQL | ug/L |
| | Chrysene | <9.6 | JU | Q | < EQL | | | | | ug/L |
| | Cobalt, total recoverable | <20 | U | | < EQL | <20 | U | | < EQL | ug/L |
| | Copper, total recoverable | <20 | U | | < EQL | <20 | U | | < EQL | ug/L |
| | Cyanide | 10 | R | Q | Rejected | | | | | ug/L |
| | Di-n-butyl phthalate | <9.6 | JU | Q | < EQL | <9.5 | JU | Q | < EQL | ug/L |
| | Di-n-octyl phthalate | <9.6 | JU | Q | < EQL | | | | | ug/L |
| | Diallate | <9.6 | JU | Q | < EQL | | | | | ug/L |
| | Dibenz(a,h)anthracene | <9.6 | JU | Q | < EQL | | | | | ug/L |
| | Dibenzofuran | <9.6 | JU | Q | < EQL | | | | | ug/L |
| | Dibromochloromethane | <25 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Dibromomethane (Methylene bromide) | <25 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Dichlorodifluoromethane | <25 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Dichloromethane (Methylene chloride) | 10 | J | I | NDD | <10 | U | | < EQL | ug/L |
| | Dieldrin | <.21 | U | | < EQL | | | | | ug/L |
| | Diethyl phthalate | <9.6 | JU | Q | < EQL | | | | | ug/L |
| | Dimethoate | <9.6 | JU | Q | < EQL | | | | | ug/L |
| | Dimethyl phthalate | <9.6 | JU | Q | < EQL | | | | | ug/L |
| | Diphenylamine | <9.6 | JU | Q | < EQL | | | | | ug/L |
| | Disulfoton | <9.6 | JU | Q | < EQL | | | | | ug/L |
| | Endosulfan I | <.1 | U | | < EQL | | | | | ug/L |
| | Endosulfan II | <.21 | U | | < EQL | | | | | ug/L |
| | Endosulfan sulfate | <.21 | U | | < EQL | | | | | ug/L |
| | Endrin | <.21 | U | | < EQL | | | | | ug/L |

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+ = exceeded the Groundwater Protection Standards listed in Appendix A.

Table B. Assessment Monitoring Results for Groundwater Wells (Cont.)
WELL LFW 34

| ANALYTICAL DATA | | | | | | |
|---------------------------------|---------------------------------|-----------------|------------|------------|-------------|-------------|
| Groundwater Protection Standard | | | | | | |
| <u>SI</u> | <u>Parameter</u> | <u>1st Half</u> | <u>CLP</u> | <u>EPA</u> | <u>File</u> | <u>Unit</u> |
| | Endrin aldehyde | <.21 | U | | < EQL | ug/L |
| | Ethyl methacrylate | <25 | U | | < EQL | ug/L |
| | Ethyl methanesulfonate | <9.6 | JU | Q | < EQL | ug/L |
| | Ethylbenzene | <25 | U | | <5 | ug/L |
| | Famphur | <200 | JU | Q | < EQL | ug/L |
| | Fluoranthene | <9.6 | JU | Q | < EQL | ug/L |
| | Fluorene | <9.6 | JU | Q | < EQL | ug/L |
| | Hepachlor | <.1 | U | | < EQL | ug/L |
| | Hepachlor epoxide | <.1 | U | | < EQL | ug/L |
| | Hexachlorobenzene | <9.6 | JU | Q | < EQL | ug/L |
| | Hexachlorobutadiene | <9.6 | JU | Q | < EQL | ug/L |
| | Hexachlorocyclopentadiene | <9.6 | JU | Q | < EQL | ug/L |
| | Hexachloroethane | <9.6 | JU | Q | < EQL | ug/L |
| | Hexachloropropene | <48 | JU | Q | < EQL | ug/L |
| | Indeno(1,2,3-c,d)pyrene | <9.6 | JU | Q | < EQL | ug/L |
| | Iodomethane (Methyl iodide) | <25 | U | | <5 | ug/L |
| | Isobutyl alcohol | <2500 | U | | < EQL | ug/L |
| | Isodrin | <9.6 | JU | Q | < EQL | ug/L |
| | Isophorone | <9.6 | JU | Q | < EQL | ug/L |
| | Isoctafrole | <9.6 | JU | Q | < EQL | ug/L |
| | Kepon | <9.6 | JU | Q | < EQL | ug/L |
| | Lead, total recoverable | <2.56 | JU | | 7.18 | ug/L |
| | Lindane | <.1 | U | | J | ug/L |
| | Mercury, total recoverable | 3.49 | U | | I | ug/L |
| | Methacrylonitrile | <1000 | U | | <5 | ug/L |
| | Methapyrene | <9.6 | U | Q | <200 | ug/L |
| | Methoxychlor | <1 | U | | < EQL | ug/L |
| | Methyl ethyl ketone | <100 | U | | < EQL | ug/L |
| | Methyl isobutyl ketone | <50 | U | | <20 | ug/L |
| | Methyl methacrylate | <100 | U | | <10 | ug/L |
| | Methyl methanesulfonate | <9.6 | JU | Q | < EQL | ug/L |
| | N-Nitrosodi-n-butylamine | <9.6 | JU | Q | < EQL | ug/L |
| | N-Nitrosodimethylamine | <9.6 | JU | Q | < EQL | ug/L |
| | N-Nitrosodiphenylamine | <24 | JU | Q | < EQL | ug/L |
| | N-Nitrosodipropylamine | <9.6 | JU | Q | < EQL | ug/L |
| | N-Nitrosomethylethylamine | <9.6 | JU | Q | < EQL | ug/L |
| | N-Nitrosopiperidine | <9.6 | JU | Q | < EQL | ug/L |
| | N-Nitrosopyrrolidine | <9.6 | JU | Q | < EQL | ug/L |
| | Naphthalene | <9.6 | JU | Q | < EQL | ug/L |
| | Nickel, total recoverable | <6.47 | JU | Q | <50 | ug/L |
| | Nitrobenzene | <9.6 | JU | Q | < EQL | ug/L |
| | O,O,O-Triethyl phosphorothioate | <9.6 | JU | Q | < EQL | ug/L |
| | PCB | | | | | |
| | PCB 1016 | <1 | U | | < EQL | ug/L |
| | PCB 1232 | <1 | U | | < EQL | ug/L |
| | PCB 1242 | <2.1 | U | | < EQL | ug/L |
| | PCB 1248 | <1 | U | | < EQL | ug/L |
| | PCB 1254 | <1 | U | | < EQL | ug/L |
| | PCB 1260 | <1 | U | | < EQL | ug/L |
| | Parathion | <1 | U | | < EQL | ug/L |
| | Parathion methyl | <9.6 | JU | Q | < EQL | ug/L |
| | Pentachlorobenzene | <9.6 | JU | Q | < EQL | ug/L |
| | Pentachloronitrobenzene | <48 | JU | Q | < EQL | ug/L |
| | Pentachlorophenol | <9.6 | JU | Q | < EQL | ug/L |
| | Phenacetin | <24 | JU | Q | < EQL | ug/L |
| | Phenanthrene | <9.6 | JU | Q | < EQL | ug/L |
| | Phenol | <9.6 | JU | Q | < EQL | ug/L |
| | Phorate | <9.6 | JU | Q | < EQL | ug/L |
| | Pronamid | <9.6 | JU | Q | < EQL | ug/L |
| | Propionitrile | <1000 | U | | < EQL | ug/L |
| | Pyrene | <9.6 | JU | Q | < EQL | ug/L |
| | Safrole | <9.6 | JU | Q | < EQL | ug/L |
| | Selenium, total recoverable | <10 | U | | <10 | ug/L |
| | Silver, total recoverable | <20 | U | | <20 | ug/L |
| | Specific conductance | 26.9 | U | | 34.9 | uS/cm |
| | Styrene | <25 | U | | <37 | ug/L |
| | Sulfide | 1000 | R | Q | < EQL | ug/L |
| | | | | | Reflected | |

Notes: Concentrations in bold italics exceed the groundwater protection standards listed in Appendix A. Bold italics were not assigned if the result was qualified with a "J", "R", "I", or "U" qualifier.
+ = exceeded the Groundwater Protection Standards listed in Appendix A.

Table B. Assessment Monitoring Results for Groundwater Wells (Cont.)

WELL LFW 34

ANALYTICAL DATA

Groundwater Protection Standard

| ST | Parameter | 1st Half | CLP | EPA | Filt. | 2nd Half | CLP | EPA | Filt. | Unit |
|----|-----------------------------|------------|-----|-----|-------|------------|-----|-----|-------|------|
| | Tetrachloroethylene | <25 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Thallium, total recoverable | <3.34 | JU | | < EQL | <10 | U | | < EQL | ug/L |
| | Thionazin | <9.6 | JU | Q | < EQL | | | | | ug/L |
| | Tin, total recoverable | <200 | U | | < EQL | | | | | ug/L |
| | Toluene | <25 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Toxaphene | <2.1 | U | | < EQL | | | | | ug/L |
| | Trichloroethylene | 2.9 | J | I | NDD | 4.7 | J | I | NDD | ug/L |
| + | Trichlorofluoromethane | 490 | | L | > 20 | 450 | | L | > 20 | ug/L |
| | Vanadium, total recoverable | <10 | U | | < EQL | <10 | U | | < EQL | ug/L |
| | Vinyl acetate | <25 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Xylenes | <50 | U | | < EQL | <10 | U | | < EQL | ug/L |
| | Zinc, total recoverable | <20 | U | | < EQL | <5.44 | JU | | < EQL | ug/L |
| | alpha-Benzene hexachloride | <.1 | U | | < EQL | | | | | ug/L |
| | alpha-Chlordane | <.1 | U | | < EQL | | | | | ug/L |
| | beta-Benzene hexachloride | <.1 | U | | < EQL | <.094 | JU | Q | < EQL | ug/L |
| | cis-1,2-Dichloroethylene | <25 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | cis-1,3-Dichloropropene | <25 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | delta-Benzene hexachloride | <.1 | U | | < EQL | | | | | ug/L |
| | gamma-Chlordane | <.1 | U | | < EQL | | | | | ug/L |
| | m-Cresol (3-Methylphenol) | | | | | | | | | |
| | m-Nitroaniline | <24 | JU | Q | < EQL | | | | | ug/L |
| | o-Cresol (2-Methylphenol) | <9.6 | JU | Q | < EQL | | | | | ug/L |
| | o-Nitroaniline | <24 | JU | Q | < EQL | | | | | ug/L |
| | o-Toluidine | <9.6 | JU | Q | < EQL | | | | | ug/L |
| | p,p'-DDD | <.21 | U | | < EQL | | | | | ug/L |
| | p,p'-DDE | <.21 | U | | < EQL | | | | | ug/L |
| | p,p'-DDT | <.21 | U | | < EQL | | | | | ug/L |
| | p-Cresol (4-Methylphenol) | <10 | JU | Q | < EQL | | | | | ug/L |
| | p-Dimethylaminoazobenzene | <9.6 | JU | Q | < EQL | | | | | ug/L |
| | p-Nitroaniline | <9.6 | JU | Q | < EQL | | | | | ug/L |
| | p-Phenylenediamine | <9.6 | JU | Q | < EQL | | | | | ug/L |
| | pH | 5.67 | J | Q | NDD | 4.94 | J | Q | NDD | pH |
| | trans-1,2-Dichloroethylene | <25 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | trans-1,3-Dichloropropene | <25 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | trans-1,4-Dichloro-2-butene | <100 | U | | < EQL | <20 | U | | < EQL | ug/L |

Notes: Concentrations in bold italics exceed the groundwater protection standards listed in Appendix A. Bold italics were not assigned if the result was qualified with a "J", "R", "L", or "U" qualifier.

+ = exceeded the Groundwater Protection Standards listed in Appendix A.

Table B. Assessment Monitoring Results for Groundwater Wells (Cont.)

WELL LFW 43B

| <u>SRS Coord.</u> | <u>Lat/Longitude</u> | <u>Screen Zone Elevation</u> | <u>Top of Casing</u> | <u>Casing</u> | <u>Pump</u> | <u>Screen Zone</u> |
|------------------------|----------------------------------|------------------------------|----------------------|---------------|-------------|--------------------|
| N 86459.2 E 45240.5 | 33.29080 Deg N -81.7156 Deg W | 100.4 - 90.4 ft msl | 203 ft msl | 4 " PVC | S | Middle |

SAMPLE DATE

02/28/00

10/02/00

FIELD DATA

| <u>Parameter</u> | <u>1st Half</u> | <u>2nd Half</u> | <u>Unit</u> |
|---------------------|-----------------|-----------------|-------------|
| Water Elevation | 163.7 | 162.8 | ft msl |
| pH | 4.4 | 5 | pH |
| Sp. Conductance | 17 | 18 | uS/cm |
| Water temperature | 18.3 | 19.3 | deg. C |
| Alkalinity as CaCO3 | 1 | 0 | mg/L |
| Turbidity | .7 | .9 | NTU |
| Volumes purged | 3.25 | 3.48 | well volume |
| Sampling code | | | |

ANALYTICAL DATA

Groundwater Protection Standard

| <u>ST</u> | <u>Parameter</u> | <u>1st Half</u> | <u>CLP</u> | <u>EPA</u> | <u>Filt.</u> | <u>2nd Half</u> | <u>CLP</u> | <u>EPA</u> | <u>Filt.</u> | <u>Unit</u> |
|-----------|--------------------------------|-----------------|------------|------------|--------------|-----------------|------------|------------|--------------|-------------|
| | 1,1,1,2-Tetrachloroethane | <5 | JU | L | < EQL | <5 | U | | < EQL | ug/L |
| | 1,1,1-Trichloroethane | <5 | JU | L | < EQL | <5 | U | | < EQL | ug/L |
| | 1,1,2,2-Tetrachloroethane | <5 | JU | L | < EQL | <5 | U | | < EQL | ug/L |
| | 1,1,2-Trichloroethane | <5 | JU | L | < EQL | <5 | U | | < EQL | ug/L |
| | 1,1-Dichloroethane | <5 | JU | L | < EQL | <5 | U | | < EQL | ug/L |
| | 1,1-Dichloroethylene | <5 | JU | L | < EQL | <5 | U | | < EQL | ug/L |
| | 1,1-Dichloropropene | <5 | JU | L | < EQL | <5 | U | | < EQL | ug/L |
| | 1,2,3-Trichloropropane | <5 | JU | L | < EQL | <5 | U | | < EQL | ug/L |
| | 1,2,4,5-Tetrachlorobenzene | <52 | U | | < EQL | | | | | ug/L |
| | 1,2,4-Trichlorobenzene | <11 | U | | < EQL | | | | | ug/L |
| | 1,2-Dibromo-3-chloropropane | <10 | JU | L | < EQL | <10 | U | | < EQL | ug/L |
| | 1,2-Dibromoethane | <5 | JU | L | < EQL | <5 | U | | < EQL | ug/L |
| | 1,2-Dichlorobenzene | <11 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,2-Dichloroethane | <5 | JU | L | < EQL | <5 | U | | < EQL | ug/L |
| | 1,2-Dichloropropane | <5 | JU | L | < EQL | <5 | U | | < EQL | ug/L |
| | 1,3,5-Trinitrobenzene | <10 | U | | < EQL | | | | | ug/L |
| | 1,3-Dichlorobenzene | <11 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,3-Dichloropropane | <5 | JU | L | < EQL | <5 | U | | < EQL | ug/L |
| | 1,3-Dinitrobenzene | <50 | U | | < EQL | | | | | ug/L |
| | 1,4-Dichlorobenzene | <11 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,4-Naphthoquinone | <11 | U | | < EQL | | | | | ug/L |
| | 1-Naphthylamine | <11 | U | | < EQL | | | | | ug/L |
| | 2,2-Dichloropropane | <5 | JU | L | < EQL | <5 | U | | < EQL | ug/L |
| | 2,2-Oxybis(1-chloropropane) | <11 | U | | < EQL | | | | | ug/L |
| | 2,3,4,6-Tetrachlorophenol | <11 | U | | < EQL | | | | | ug/L |
| | 2,4,5-T | | | | | | | | | |
| | 2,4,5-TP (Silvex) | | | | | | | | | |
| | 2,4,5-Trichlorophenol | <11 | U | | < EQL | | | | | ug/L |
| | 2,4,6-Trichlorophenol | <26 | U | | < EQL | | | | | ug/L |
| | 2,4-Dichlorophenol | <11 | U | | < EQL | | | | | ug/L |
| | 2,4-Dichlorophenoxyacetic acid | | | | | | | | | |
| | 2,4-Dimethyl phenol | <11 | U | | < EQL | | | | | ug/L |
| | 2,4-Dinitrophenol | <26 | U | | < EQL | | | | | ug/L |
| | 2,4-Dinitrotoluene | <11 | U | | < EQL | | | | | ug/L |
| | 2,6-Dichlorophenol | <25 | U | | < EQL | | | | | ug/L |
| | 2,6-Dinitrotoluene | <11 | U | | < EQL | | | | | ug/L |
| | 2-Acetylaminofluorene | <11 | U | | < EQL | | | | | ug/L |
| | 2-Chloronaphthalene | <11 | U | | < EQL | | | | | ug/L |
| | 2-Chlorophenol | <11 | U | | < EQL | | | | | ug/L |
| | 2-Hexanone | <20 | JU | L | < EQL | <20 | U | | < EQL | ug/L |
| | 2-Methyl-4,6-dinitrophenol | <26 | U | | < EQL | | | | | ug/L |
| | 2-Methylnaphthalene | <11 | U | | < EQL | | | | | ug/L |
| | 2-Naphthylamine | <11 | U | | < EQL | | | | | ug/L |
| | 2-Nitrophenol | <11 | U | | < EQL | | | | | ug/L |
| | 2-sec-Butyl-4,6-dinitrophenol | | | | | | | | | |
| | 3,3"-Dichlorobenzidine | <11 | U | | < EQL | | | | | ug/L |
| | 3,3"-Dimethylbenzidine | <21 | U | | < EQL | | | | | ug/L |
| | 3-Methylcholanthrene | <11 | U | | < EQL | | | | | ug/L |
| | 4-Aminobiphenyl | <11 | U | | < EQL | | | | | ug/L |
| | 4-Bromophenyl phenyl ether | <11 | U | | < EQL | | | | | ug/L |
| | 4-Chloro-m-cresol | <11 | U | | < EQL | | | | | ug/L |

Notes: Concentrations in bold italics exceed the groundwater protection standards listed in Appendix A. Bold italics were not assigned if the result was qualified with a "J", "R", "L", or "U" qualifier.

+ = exceeded the Groundwater Protection Standards listed in Appendix A.

Table B. Assessment Monitoring Results for Groundwater Wells (Cont.)

WELL LFW 43B

ANALYTICAL DATA

Groundwater Protection Standard

| ST | Parameter | 1st Half | CLP | EPA | Filt. | 2nd Half | CLP | EPA | Filt. | Unit |
|----|--------------------------------------|----------|-----|-----|-------|----------|-----|-----|--------|------|
| | 4-Chloroaniline | <11 | U | | < EQL | | | | | ug/L |
| | 4-Chlorophenyl phenyl ether | <11 | U | | < EQL | | | | | ug/L |
| | 4-Nitrophenol | <26 | U | | < EQL | | | | | ug/L |
| | 5-Nitro-o-toluidine | <11 | U | | < EQL | | | | | ug/L |
| | 7,12-Dimethylbenz(a)anthracene | <11 | U | | < EQL | | | | | ug/L |
| | Acenaphthene | <11 | U | | < EQL | | | | | ug/L |
| | Acenaphthylene | <11 | U | | < EQL | | | | | ug/L |
| | Acetone | <20 | JU | L | < EQL | <20 | U | | < EQL | ug/L |
| | Acetonitrile (Methyl cyanide) | <200 | JU | L | < EQL | <200 | U | | < EQL | ug/L |
| | Acetophenone | <11 | U | | < EQL | | | | | ug/L |
| | Acrolein | <50 | JU | L | < EQL | <50 | U | | < EQL | ug/L |
| | Acrylonitrile | <10 | JU | L | < EQL | <10 | U | | < EQL | ug/L |
| | Aldrin | <.1 | U | | < EQL | | | | | ug/L |
| | Allyl chloride | <5 | JU | L | < EQL | <10 | U | | < EQL | ug/L |
| | Anthracene | <11 | U | | < EQL | | | | | ug/L |
| | Antimony, total recoverable | <100 | U | | < EQL | <100 | U | | < EQL | ug/L |
| | Arsenic, total recoverable | <5.01 | JU | | < EQL | <40 | U | | < EQL | ug/L |
| | Barium, total recoverable | 3.52 | J | I | NDD | 4 | | | < 2000 | ug/L |
| | Benzene | <5 | JU | L | < EQL | <5 | U | | < EQL | ug/L |
| | Benzo(a)anthracene | <11 | U | | < EQL | | | | | ug/L |
| | Benzo(a)pyrene | <11 | U | | < EQL | | | | | ug/L |
| | Benzo(b)fluoranthene | <11 | U | | < EQL | | | | | ug/L |
| | Benzo(g,h,i)perylene | <11 | U | | < EQL | | | | | ug/L |
| | Benzo(k)fluoranthene | <11 | U | | < EQL | | | | | ug/L |
| | Benzyl alcohol | | | | | | | | | |
| | Beryllium, total recoverable | <1 | U | | < EQL | <.13 | JU | | < EQL | ug/L |
| | Bis(2-chloroethoxy) methane | <11 | U | | < EQL | | | | | ug/L |
| | Bis(2-chloroethyl) ether | <11 | U | | < EQL | | | | | ug/L |
| | Bis(2-ethylhexyl) phthalate | <11 | U | | < EQL | | | | | ug/L |
| | Bromochloromethane | <5 | JU | L | < EQL | <5 | U | | < EQL | ug/L |
| | Bromodichloromethane | <5 | JU | L | < EQL | <5 | U | | < EQL | ug/L |
| | Bromoform | <5 | JU | L | < EQL | <5 | U | | < EQL | ug/L |
| | Bromomethane (Methyl bromide) | <5 | JU | L | < EQL | <10 | U | | < EQL | ug/L |
| | Butylbenzyl phthalate | <11 | U | | < EQL | | | | | ug/L |
| | Cadmium, total recoverable | <10 | U | | < EQL | <10 | U | | < EQL | ug/L |
| | Carbon disulfide | <5 | JU | L | < EQL | <5 | U | | < EQL | ug/L |
| | Carbon tetrachloride | <5 | JU | L | < EQL | <5 | U | | < EQL | ug/L |
| | Chlorobenzene | <5 | JU | L | < EQL | <5 | U | | < EQL | ug/L |
| | Chlorobenzilate | <11 | U | | < EQL | | | | | ug/L |
| | Chloroethane | <5 | JU | L | < EQL | <10 | U | | < EQL | ug/L |
| | Chloroethene (Vinyl chloride) | <5 | JU | L | < EQL | <10 | U | | < EQL | ug/L |
| | Chloroform | <5 | JU | L | < EQL | <5 | U | | < EQL | ug/L |
| | Chloromethane (Methyl chloride) | <5 | JU | L | < EQL | <10 | U | | < EQL | ug/L |
| | Chloroprene | <20 | JU | L | < EQL | <20 | U | | < EQL | ug/L |
| | Chromium, total recoverable | <10 | U | | < EQL | <10 | U | | < EQL | ug/L |
| | Chrysene | <11 | U | | < EQL | | | | | ug/L |
| | Cobalt, total recoverable | <6.25 | JU | V | < EQL | <20 | U | | < EQL | ug/L |
| | Copper, total recoverable | <2.83 | JU | V | < EQL | <20 | U | | < EQL | ug/L |
| | Cyanide | | | | | | | | | |
| | Di-n-butyl phthalate | <11 | U | | < EQL | <9.5 | U | | < EQL | ug/L |
| | Di-n-octyl phthalate | <11 | U | | < EQL | | | | | ug/L |
| | Diallate | <11 | U | | < EQL | | | | | ug/L |
| | Dibenz(a,h)anthracene | <11 | U | | < EQL | | | | | ug/L |
| | Dibenzofuran | <11 | U | | < EQL | | | | | ug/L |
| | Dibromochloromethane | <5 | JU | L | < EQL | <5 | U | | < EQL | ug/L |
| | Dibromomethane (Methylene bromide) | <5 | JU | L | < EQL | <5 | U | | < EQL | ug/L |
| | Dichlorodifluoromethane | <5 | JU | L | < EQL | <10 | U | | < EQL | ug/L |
| + | Dichloromethane (Methylene chloride) | <2.8 | JU | LV | < EQL | 8.32 | | | > 5 | ug/L |
| | Dieldrin | <.21 | U | | < EQL | | | | | ug/L |
| | Diethyl phthalate | <11 | U | | < EQL | | | | | ug/L |
| | Dimethoate | <11 | U | | < EQL | | | | | ug/L |
| | Dimethyl phthalate | <11 | U | | < EQL | | | | | ug/L |
| | Diphenylamine | <11 | U | | < EQL | | | | | ug/L |
| | Disulfoton | <11 | U | | < EQL | | | | | ug/L |
| | Endosulfan I | <.1 | U | | < EQL | | | | | ug/L |
| | Endosulfan II | <.21 | U | | < EQL | | | | | ug/L |
| | Endosulfan sulfate | <.21 | U | | < EQL | | | | | ug/L |
| | Endrin | <.21 | U | | < EQL | | | | | ug/L |

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+ = exceeded the Groundwater Protection Standards listed in Appendix A.

Table B. Assessment Monitoring Results for Groundwater Wells (Cont.)

WELL LFW 43B

ANALYTICAL DATA

Groundwater Protection Standard

| ST | Parameter | 1st Half | CLP | EPA | Filt. | 2nd Half | CLP | EPA | Filt. | Unit |
|----|---------------------------------|----------|-----|-----|-------|----------|-----|-----|-------|-------|
| | Endrin aldehyde | <.21 | U | | < EQL | | | | | ug/L |
| | Ethyl methacrylate | <5 | JU | L | < EQL | <5 | U | | < EQL | ug/L |
| | Ethyl methanesulfonate | <11 | U | | < EQL | | | | | ug/L |
| | Ethylbenzene | <5 | JU | L | < EQL | <5 | U | | < EQL | ug/L |
| | Famphur | <200 | U | | < EQL | | | | | ug/L |
| | Fluoranthene | <11 | U | | < EQL | | | | | ug/L |
| | Fluorene | <11 | U | | < EQL | | | | | ug/L |
| | Heptachlor | <.1 | U | | < EQL | | | | | ug/L |
| | Heptachlor epoxide | <.1 | U | | < EQL | | | | | ug/L |
| | Hexachlorobenzene | <11 | U | | < EQL | | | | | ug/L |
| | Hexachlorobutadiene | <11 | U | | < EQL | | | | | ug/L |
| | Hexachlorocyclopentadiene | <11 | U | | < EQL | | | | | ug/L |
| | Hexachloroethane | <11 | U | | < EQL | | | | | ug/L |
| | Hexachloropropene | <52 | U | | < EQL | | | | | ug/L |
| | Indeno(1,2,3-c,d)pyrene | <11 | U | | < EQL | | | | | ug/L |
| | Iodomethane (Methyl iodide) | <5 | JU | L | < EQL | <5 | U | | < EQL | ug/L |
| | Isobutyl alcohol | <500 | JU | L | < EQL | <500 | U | | < EQL | ug/L |
| | Isodrin | <11 | U | | < EQL | | | | | ug/L |
| | Isophorone | <11 | U | | < EQL | | | | | ug/L |
| | Isosafrole | <11 | U | | < EQL | | | | | ug/L |
| | Kepone | <11 | U | | < EQL | | | | | ug/L |
| | Lead, total recoverable | <10 | U | | < EQL | <47 | U | | < EQL | ug/L |
| | Lindane | <.1 | U | | < EQL | | | | | ug/L |
| | Mercury, total recoverable | <.5 | U | | < EQL | <.5 | U | | < EQL | ug/L |
| | Methacrylonitrile | <200 | JU | L | < EQL | <200 | U | | < EQL | ug/L |
| | Methapyrilene | <11 | U | | < EQL | | | | | ug/L |
| | Methoxychlor | <1 | U | | < EQL | | | | | ug/L |
| | Methyl ethyl ketone | <20 | JU | L | < EQL | <20 | U | | < EQL | ug/L |
| | Methyl isobutyl ketone | <10 | JU | L | < EQL | <10 | U | | < EQL | ug/L |
| | Methyl methacrylate | <20 | JU | L | < EQL | <20 | U | | < EQL | ug/L |
| | Methyl methanesulfonate | <11 | U | | < EQL | | | | | ug/L |
| | N-Nitrosodi-n-butylamine | <11 | U | | < EQL | | | | | ug/L |
| | N-Nitrosodiethylamine | <11 | U | | < EQL | | | | | ug/L |
| | N-Nitrosodimethylamine | <26 | U | | < EQL | | | | | ug/L |
| | N-Nitrosodiphenylamine | <11 | U | | < EQL | | | | | ug/L |
| | N-Nitrosodipropylamine | <11 | U | | < EQL | | | | | ug/L |
| | N-Nitrosomethylethylamine | <11 | U | | < EQL | | | | | ug/L |
| | N-Nitrosopiperidine | <11 | U | | < EQL | | | | | ug/L |
| | N-Nitrosopyrrolidine | <11 | U | | < EQL | | | | | ug/L |
| | Naphthalene | <11 | U | | < EQL | | | | | ug/L |
| | Nickel, total recoverable | <50 | U | | < EQL | <50 | U | | < EQL | ug/L |
| | Nitrobenzene | <11 | U | | < EQL | | | | | ug/L |
| | O,O,O-Triethyl phosphorothioate | <11 | U | | < EQL | | | | | ug/L |
| | PCB | | | | | | | | | |
| | PCB 1016 | | | | | | | | | |
| | PCB 1232 | | | | | | | | | |
| | PCB 1242 | | | | | | | | | |
| | PCB 1248 | | | | | | | | | |
| | PCB 1254 | | | | | | | | | |
| | PCB 1260 | | | | | | | | | |
| | Parathion | <11 | U | | < EQL | | | | | ug/L |
| | Parathion methyl | <11 | U | | < EQL | | | | | ug/L |
| | Pentachlorobenzene | <52 | U | | < EQL | | | | | ug/L |
| | Pentachloronitrobenzene | <11 | U | | < EQL | | | | | ug/L |
| | Pentachlorophenol | <26 | U | | < EQL | | | | | ug/L |
| | Phenacetin | <11 | U | | < EQL | | | | | ug/L |
| | Phenanthrene | <11 | U | | < EQL | | | | | ug/L |
| | Phenol | <11 | U | | < EQL | | | | | ug/L |
| | Phorate | <11 | U | | < EQL | | | | | ug/L |
| | Pronamid | <11 | U | | < EQL | | | | | ug/L |
| | Propionitrile | <200 | JU | L | < EQL | <200 | U | | < EQL | ug/L |
| | Pyrene | <11 | U | | < EQL | | | | | ug/L |
| | Safrole | <11 | U | | < EQL | | | | | ug/L |
| | Selenium, total recoverable | <10 | U | | < EQL | <66 | U | | < EQL | ug/L |
| | Silver, total recoverable | <20 | U | | < EQL | 3.59 | J | I | NDD | ug/L |
| | Specific conductance | 16.1 | | | < 37 | 16.1 | | | < 37 | uS/cm |
| | Styrene | <5 | JU | L | < EQL | <5 | U | | < EQL | ug/L |
| | Sulfide | | | | | | | | | |

Notes: Concentrations in bold italics exceed the groundwater protection standards listed in Appendix A. Bold italics were not assigned if the result was qualified with a "J", "R", "L", or "U" qualifier.

+ = exceeded the Groundwater Protection Standards listed in Appendix A.

Table B. Assessment Monitoring Results for Groundwater Wells (Cont.)

WELL LFW 43B

ANALYTICAL DATA

Groundwater Protection Standard

| ST | Parameter | 1st Half | CLP | EPA | Filt. | 2nd Half | CLP | EPA | Filt. | Unit |
|-----------|-----------------------------|-----------------|------------|------------|--------------|-----------------|------------|------------|--------------|-------------|
| | Tetrachloroethylene | <5 | JU | L | < EQL | <5 | U | | < EQL | ug/L |
| | Thallium, total recoverable | <10 | U | | < EQL | <10 | U | | < EQL | ug/L |
| | Thionazin | <11 | U | | < EQL | | | | | ug/L |
| | Tin, total recoverable | <200 | U | | < EQL | | | | | ug/L |
| | Toluene | <5 | JU | L | < EQL | <5 | U | | < EQL | ug/L |
| | Toxaphene | <2.1 | U | | < EQL | | | | | ug/L |
| | Trichloroethylene | <5 | JU | L | < EQL | <5 | U | | < EQL | ug/L |
| | Trichlorofluoromethane | <5 | JU | L | < EQL | <5 | U | | < EQL | ug/L |
| | Vanadium, total recoverable | <10 | U | | < EQL | <10 | U | | < EQL | ug/L |
| | Vinyl acetate | <5 | JU | L | < EQL | <10 | U | | < EQL | ug/L |
| | Xylenes | <10 | JU | L | < EQL | <10 | U | | < EQL | ug/L |
| | Zinc, total recoverable | 10.2 | J | I | NDD | <20 | U | | < EQL | ug/L |
| | alpha-Benzene hexachloride | <.1 | U | | < EQL | | | | | ug/L |
| | alpha-Chlordane | <.1 | U | | < EQL | | | | | ug/L |
| | beta-Benzene hexachloride | <.1 | U | | < EQL | <.095 | U | | < EQL | ug/L |
| | cis-1,2-Dichloroethylene | <5 | JU | L | < EQL | <5 | U | | < EQL | ug/L |
| | cis-1,3-Dichloropropene | <5 | JU | L | < EQL | <5 | U | | < EQL | ug/L |
| | delta-Benzene hexachloride | <.1 | U | | < EQL | | | | | ug/L |
| | gamma-Chlordane | <.1 | U | | < EQL | | | | | ug/L |
| | m-Cresol (3-Methylphenol) | | | | | | | | | |
| | m-Nitroaniline | <26 | U | | < EQL | | | | | ug/L |
| | o-Cresol (2-Methylphenol) | <11 | U | | < EQL | | | | | ug/L |
| | o-Nitroaniline | <26 | U | | < EQL | | | | | ug/L |
| | o-Toluidine | <11 | U | | < EQL | | | | | ug/L |
| | p,p"-DDD | <.21 | U | | < EQL | | | | | ug/L |
| | p,p"-DDE | <.21 | U | | < EQL | | | | | ug/L |
| | p,p"-DDT | <.21 | U | | < EQL | | | | | ug/L |
| | p-Cresol (4-Methylphenol) | <10 | U | | < EQL | | | | | ug/L |
| | p-Dimethylaminoazobenzene | <11 | U | | < EQL | | | | | ug/L |
| | p-Nitroaniline | <11 | U | | < EQL | | | | | ug/L |
| | p-Phenylenediamine | <11 | U | | < EQL | | | | | ug/L |
| | pH | 5.64 | J | Q | NDD | 5.04 | J | Q | NDD | pH |
| | trans-1,2-Dichloroethylene | <5 | JU | L | < EQL | <5 | U | | < EQL | ug/L |
| | trans-1,3-Dichloropropene | <5 | JU | L | < EQL | <5 | U | | < EQL | ug/L |
| | trans-1,4-Dichloro-2-butene | <20 | JU | L | < EQL | <20 | U | | < EQL | ug/L |

Notes: Concentrations in bold italics exceed the groundwater protection standards listed in Appendix A. Bold italics were not assigned if the result was qualified with a "J", "R", "L", or "U" qualifier.

+ = exceeded the Groundwater Protection Standards listed in Appendix A.

Table B. Assessment Monitoring Results for Groundwater Wells (Cont.)

WELL LFW 43C

| <u>SRS Coord.</u> | <u>Lat/Longitude</u> | <u>Screen Zone Elevation</u> | <u>Top of Casing</u> | <u>Casing</u> | <u>Pump</u> | <u>Screen Zone</u> |
|------------------------|----------------------------------|------------------------------|----------------------|---------------|-------------|--------------------|
| N 86480.6 E 45234.9 | 33.29084 Deg N -81.7157 Deg W | 138.5 - 128.5 ft msl | 202.6 ft msl | 4 " PVC | S | Upper |

SAMPLE DATE

02/28/00

10/02/00

FIELD DATA

| <u>Parameter</u> | <u>1st Half</u> | <u>2nd Half</u> | <u>Unit</u> |
|---------------------|-----------------|-----------------|-------------|
| Water Elevation | 163.69 | 162.83 | ft msl |
| pH | 4.4 | 5.2 | pH |
| Sp. Conductance | 14 | 15 | uS/cm |
| Water temperature | 17.6 | 19.3 | deg. C |
| Alkalinity as CaCO3 | 1 | 0 | mg/L |
| Turbidity | .7 | 1.3 | NTU |
| Volumes purged | 4.01 | 4.78 | well volume |
| Sampling code | | | |

ANALYTICAL DATA

Groundwater Protection Standard

| <u>ST</u> | <u>Parameter</u> | <u>1st Half</u> | <u>CLP</u> | <u>EPA</u> | <u>Filt.</u> | <u>2nd Half</u> | <u>CLP</u> | <u>EPA</u> | <u>Filt.</u> | <u>Unit</u> |
|-----------|--------------------------------|-----------------|------------|------------|--------------|-----------------|------------|------------|--------------|-------------|
| | 1,1,1,2-Tetrachloroethane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,1,1-Trichloroethane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,1,2,2-Tetrachloroethane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,1,2-Trichloroethane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,1-Dichloroethane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,1-Dichloroethylene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,1-Dichloropropene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,2,3-Trichloropropane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,2,4,5-Tetrachlorobenzene | <52 | U | | < EQL | | | | | ug/L |
| | 1,2,4-Trichlorobenzene | <11 | U | | < EQL | | | | | ug/L |
| | 1,2-Dibromo-3-chloropropane | <10 | U | | < EQL | <10 | U | | < EQL | ug/L |
| | 1,2-Dibromoethane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,2-Dichlorobenzene | <11 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,2-Dichloroethane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,2-Dichloropropane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,3,5-Trinitrobenzene | <10 | U | | < EQL | | | | | ug/L |
| | 1,3-Dichlorobenzene | <11 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,3-Dichloropropane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,3-Dinitrobenzene | <50 | U | | < EQL | | | | | ug/L |
| | 1,4-Dichlorobenzene | <11 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,4-Naphthoquinone | <11 | U | | < EQL | | | | | ug/L |
| | 1-Naphthylamine | <11 | U | | < EQL | | | | | ug/L |
| | 2,2-Dichloropropane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 2,2-Oxybis(1-chloropropane) | <11 | U | | < EQL | | | | | ug/L |
| | 2,3,4,6-Tetrachlorophenol | <11 | U | | < EQL | | | | | ug/L |
| | 2,4,5-T | | | | | | | | | |
| | 2,4,5-TP (Silvex) | | | | | | | | | |
| | 2,4,5-Trichlorophenol | <11 | U | | < EQL | | | | | ug/L |
| | 2,4,6-Trichlorophenol | <26 | U | | < EQL | | | | | ug/L |
| | 2,4-Dichlorophenol | <11 | U | | < EQL | | | | | ug/L |
| | 2,4-Dichlorophenoxyacetic acid | | | | | | | | | |
| | 2,4-Dimethyl phenol | <11 | U | | < EQL | | | | | ug/L |
| | 2,4-Dinitrophenol | <26 | U | | < EQL | | | | | ug/L |
| | 2,4-Dinitrotoluene | <11 | U | | < EQL | | | | | ug/L |
| | 2,6-Dichlorophenol | <25 | U | | < EQL | | | | | ug/L |
| | 2,6-Dinitrotoluene | <11 | U | | < EQL | | | | | ug/L |
| | 2-Acetylaminofluorene | <11 | U | | < EQL | | | | | ug/L |
| | 2-Chloronaphthalene | <11 | U | | < EQL | | | | | ug/L |
| | 2-Chlorophenol | <11 | U | | < EQL | | | | | ug/L |
| | 2-Hexanone | <20 | U | | < EQL | <20 | U | | < EQL | ug/L |
| | 2-Methyl-4,6-dinitrophenol | <26 | U | | < EQL | | | | | ug/L |
| | 2-Methylnaphthalene | <11 | U | | < EQL | | | | | ug/L |
| | 2-Naphthylamine | <11 | U | | < EQL | | | | | ug/L |
| | 2-Nitrophenol | <11 | U | | < EQL | | | | | ug/L |
| | 2-sec-Butyl-4,6-dinitrophenol | | | | | | | | | |
| | 3,3"-Dichlorobenzidine | <11 | U | | < EQL | | | | | ug/L |
| | 3,3"-Dimethylbenzidine | <21 | U | | < EQL | | | | | ug/L |
| | 3-Methylcholanthrene | <11 | U | | < EQL | | | | | ug/L |
| | 4-Aminobiphenyl | <11 | U | | < EQL | | | | | ug/L |
| | 4-Bromophenyl phenyl ether | <11 | U | | < EQL | | | | | ug/L |
| | 4-Chloro-m-cresol | <11 | U | | < EQL | | | | | ug/L |

Notes: Concentrations in bold italics exceed the groundwater protection standards listed in Appendix A. Bold italics were not assigned if the result was qualified with a "J", "R", "L", or "U" qualifier.

+ = exceeded the Groundwater Protection Standards listed in Appendix A.

Table B. Assessment Monitoring Results for Groundwater Wells (Cont.)

WELL LFW 43C

ANALYTICAL DATA

Groundwater Protection Standard

| ST | Parameter | 1st Half | CLP | EPA | Filt. | 2nd Half | CLP | EPA | Filt. | Unit |
|----|--------------------------------------|----------|-----|-----|-------|----------|-----|-----|-------|------|
| | 4-Chloroaniline | <11 | U | | < EQL | | | | | ug/L |
| | 4-Chlorophenyl phenyl ether | <11 | U | | < EQL | | | | | ug/L |
| | 4-Nitrophenol | <26 | U | | < EQL | | | | | ug/L |
| | 5-Nitro-o-toluidine | <11 | U | | < EQL | | | | | ug/L |
| | 7,12-Dimethylbenz(a)anthracene | <11 | U | | < EQL | | | | | ug/L |
| | Acenaphthene | <11 | U | | < EQL | | | | | ug/L |
| | Acenaphthylene | <11 | U | | < EQL | | | | | ug/L |
| | Acetone | <20 | U | | < EQL | <20 | U | | < EQL | ug/L |
| | Acetonitrile (Methyl cyanide) | <200 | U | | < EQL | <200 | U | | < EQL | ug/L |
| | Acetophenone | <11 | U | | < EQL | | | | | ug/L |
| | Acrolein | <50 | U | | < EQL | <50 | U | | < EQL | ug/L |
| | Acrylonitrile | <10 | U | | < EQL | <10 | U | | < EQL | ug/L |
| | Aldrin | <.1 | U | | < EQL | | | | | ug/L |
| | Allyl chloride | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Anthracene | <11 | U | | < EQL | | | | | ug/L |
| | Antimony, total recoverable | <100 | U | | < EQL | <100 | U | | < EQL | ug/L |
| | Arsenic, total recoverable | <10 | U | | < EQL | 4.48 | J | I | NDD | ug/L |
| | Barium, total recoverable | 2.99 | J | I | NDD | 3.42 | J | I | NDD | ug/L |
| | Benzene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Benzo(a)anthracene | <11 | U | | < EQL | | | | | ug/L |
| | Benzo(a)pyrene | <11 | U | | < EQL | | | | | ug/L |
| | Benzo(b)fluoranthene | <11 | U | | < EQL | | | | | ug/L |
| | Benzo(g,h,i)perylene | <11 | U | | < EQL | | | | | ug/L |
| | Benzo(k)fluoranthene | <11 | U | | < EQL | | | | | ug/L |
| | Benzyl alcohol | | | | | | | | | |
| | Beryllium, total recoverable | <1 | U | | < EQL | <.13 | JU | | < EQL | ug/L |
| | Bis(2-chloroethoxy) methane | <11 | U | | < EQL | | | | | ug/L |
| | Bis(2-chloroethyl) ether | <11 | U | | < EQL | | | | | ug/L |
| | Bis(2-ethylhexyl) phthalate | 6.6 | J | I | NDD | | | | | ug/L |
| | Bromochloromethane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Bromodichloromethane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Bromoform | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Bromomethane (Methyl bromide) | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Butylbenzyl phthalate | <11 | U | | < EQL | | | | | ug/L |
| | Cadmium, total recoverable | <10 | U | | < EQL | <10 | U | | < EQL | ug/L |
| | Carbon disulfide | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Carbon tetrachloride | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Chlorobenzene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Chlorobenzilate | <11 | U | | < EQL | | | | | ug/L |
| | Chloroethane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Chloroethene (Vinyl chloride) | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Chloroform | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Chloromethane (Methyl chloride) | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Chloroprene | <20 | U | | < EQL | <20 | U | | < EQL | ug/L |
| | Chromium, total recoverable | <10 | U | | < EQL | <10 | U | | < EQL | ug/L |
| | Chrysene | <11 | U | | < EQL | | | | | ug/L |
| | Cobalt, total recoverable | <5.19 | JU | V | < EQL | <5.26 | JU | | < EQL | ug/L |
| | Copper, total recoverable | <3.96 | JU | | < EQL | <20 | U | | < EQL | ug/L |
| | Cyanide | | | | | | | | | |
| | Di-n-butyl phthalate | <11 | U | | < EQL | <9.5 | U | | < EQL | ug/L |
| | Di-n-octyl phthalate | <11 | U | | < EQL | | | | | ug/L |
| | Diallylate | <11 | U | | < EQL | | | | | ug/L |
| | Dibenz(a,h)anthracene | <11 | U | | < EQL | | | | | ug/L |
| | Dibenzofuran | <11 | U | | < EQL | | | | | ug/L |
| | Dibromochloromethane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Dibromomethane (Methylene bromide) | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Dichlorodifluoromethane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Dichloromethane (Methylene chloride) | <10 | U | | < EQL | <10 | U | | < EQL | ug/L |
| | Dieldrin | <.21 | U | | < EQL | | | | | ug/L |
| | Diethyl phthalate | <11 | U | | < EQL | | | | | ug/L |
| | Dimethoate | <11 | U | | < EQL | | | | | ug/L |
| | Dimethyl phthalate | <11 | U | | < EQL | | | | | ug/L |
| | Diphenylamine | <11 | U | | < EQL | | | | | ug/L |
| | Disulfoton | <11 | U | | < EQL | | | | | ug/L |
| | Endosulfan I | <.1 | U | | < EQL | | | | | ug/L |
| | Endosulfan II | <.21 | U | | < EQL | | | | | ug/L |
| | Endosulfan sulfate | <.21 | U | | < EQL | | | | | ug/L |
| | Endrin | <.21 | U | | < EQL | | | | | ug/L |

Notes: Concentrations in bold italics exceed the groundwater protection standards listed in Appendix A. Bold italics were not assigned if the result was qualified with a "J", "R", "L", or "U" qualifier.

+ = exceeded the Groundwater Protection Standards listed in Appendix A.

Table B. Assessment Monitoring Results for Groundwater Wells (Cont.)
WELL LPW 43C

| ANALYTICAL DATA | | | | | | |
|---------------------------------|---------------------------------|----------|-----|-----|--------|-------|
| Groundwater Protection Standard | | | | | | |
| SI | Parameter | 1st Half | CLP | EPA | Filter | Unit |
| | Endrin aldehyde | <21 | U | | < EQL | ug/L |
| | Ethyl methacrylate | <5 | U | | < EQL | ug/L |
| | Ethyl methanesulfonate | <11 | U | | < EQL | ug/L |
| | Ethylbenzene | <5 | U | | < EQL | ug/L |
| | Famphur | <200 | U | | < EQL | ug/L |
| | Fluoranthene | <11 | U | | < EQL | ug/L |
| | Fluorene | <11 | U | | < EQL | ug/L |
| | Heptachlor | <1 | U | | < EQL | ug/L |
| | Heptachlor epoxide | <1 | U | | < EQL | ug/L |
| | Hexachlorobenzene | <11 | U | | < EQL | ug/L |
| | Hexachlorobutadiene | <11 | U | | < EQL | ug/L |
| | Hexachlorocyclopentadiene | <11 | U | | < EQL | ug/L |
| | Hexachloroethane | <11 | U | | < EQL | ug/L |
| | Hexachloropropene | <52 | U | | < EQL | ug/L |
| | Indeno (1,2,3-c,d)pyrene | <11 | U | | < EQL | ug/L |
| | Iodomethane (Methyl iodide) | <5 | U | | < EQL | ug/L |
| | Isobutyl alcohol | <500 | U | | < EQL | ug/L |
| | Isodrin | <11 | U | | < EQL | ug/L |
| | Isophorone | <11 | U | | < EQL | ug/L |
| | Isosafrole | <11 | U | | < EQL | ug/L |
| | Kepone | <11 | U | | < EQL | ug/L |
| | Lead, total recoverable | <10 | U | | < EQL | ug/L |
| | Lindane | <1 | U | | < EQL | ug/L |
| | Mercury, total recoverable | <5 | U | | < EQL | ug/L |
| | Methacrylonitrile | <200 | U | | < EQL | ug/L |
| | Methapyrene | <11 | U | | < EQL | ug/L |
| | Methoxychlor | <1 | U | | < EQL | ug/L |
| | Methyl ethyl ketone | <20 | U | | < EQL | ug/L |
| | Methyl isobutyl ketone | <10 | U | | < EQL | ug/L |
| | Methyl methacrylate | <20 | U | | < EQL | ug/L |
| | Methyl methanesulfonate | <11 | U | | < EQL | ug/L |
| | N-Nitrosodi-n-butylamine | <11 | U | | < EQL | ug/L |
| | N-Nitrosodimethylamine | <11 | U | | < EQL | ug/L |
| | N-Nitrosodiphenylamine | <26 | U | | < EQL | ug/L |
| | N-Nitrosodipropylamine | <11 | U | | < EQL | ug/L |
| | N-Nitrosomethylethylamine | <11 | U | | < EQL | ug/L |
| | N-Nitrosopiperidine | <11 | U | | < EQL | ug/L |
| | N-Nitrosopyrrolidine | <11 | U | | < EQL | ug/L |
| | Napthalene | <11 | U | | < EQL | ug/L |
| | Nickel, total recoverable | <50 | U | | < EQL | ug/L |
| | Nitrobenzene | <11 | U | | < EQL | ug/L |
| | O,O,O-Triethyl phosphorothioate | <11 | U | | < EQL | ug/L |
| | PCB | | | | | |
| | PCB 1016 | | | | | |
| | PCB 1232 | | | | | |
| | PCB 1242 | | | | | |
| | PCB 1248 | | | | | |
| | PCB 1254 | | | | | |
| | PCB 1260 | | | | | |
| | Parathion | <11 | U | | < EQL | ug/L |
| | Parathion methyl | <11 | U | | < EQL | ug/L |
| | Pentachlorobenzene | <52 | U | | < EQL | ug/L |
| | Pentachloronitrobenzene | <11 | U | | < EQL | ug/L |
| | Pentachlorophenol | <26 | U | | < EQL | ug/L |
| | Phenacetin | <11 | U | | < EQL | ug/L |
| | Phenanthrene | <11 | U | | < EQL | ug/L |
| | Phenol | <11 | U | | < EQL | ug/L |
| | Phorate | <11 | U | | < EQL | ug/L |
| | Pronamid | <11 | U | | < EQL | ug/L |
| | Propionitrile | <200 | U | | < EQL | ug/L |
| | Pyrene | <11 | U | | < EQL | ug/L |
| | Safrole | <11 | U | | < EQL | ug/L |
| | Selenium, total recoverable | <10 | U | | < EQL | ug/L |
| | Silver, total recoverable | <4.67 | U | | < EQL | ug/L |
| | Specific conductance | 13.4 | | | < 37 | uS/cm |
| | Styrene | <5 | U | | < EQL | ug/L |
| | Sulfide | | | | | |

Notes: Concentrations in bold italics exceed the groundwater protection standards listed in Appendix A. Bold italics were not assigned if the result was qualified with a "J", "R", "L", or "U" qualifier.
+ = exceeded the Groundwater Protection Standards listed in Appendix A.

Table B. Assessment Monitoring Results for Groundwater Wells (Cont.)

WELL LFW 43C

ANALYTICAL DATA

Groundwater Protection Standard

| ST | Parameter | 1st Half | CLP | EPA | Filt. | 2nd Half | CLP | EPA | Filt. | Unit |
|----|-----------------------------|----------|-----|-----|-------|----------|-----|-----|-------|------|
| | Tetrachloroethylene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Thallium, total recoverable | <7.36 | JU | | < EQL | 8.05 | J | I | NDD | ug/L |
| | Thionazin | <11 | U | | < EQL | | | | | ug/L |
| | Tin, total recoverable | <200 | U | | < EQL | | | | | ug/L |
| | Toluene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Toxaphene | <2.1 | U | | < EQL | | | | | ug/L |
| | Trichloroethylene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Trichlorofluoromethane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Vanadium, total recoverable | <10 | U | | < EQL | <10 | U | | < EQL | ug/L |
| | Vinyl acetate | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Xylenes | <10 | U | | < EQL | <10 | U | | < EQL | ug/L |
| | Zinc, total recoverable | <20 | U | | < EQL | <20 | U | | < EQL | ug/L |
| | alpha-Benzene hexachloride | <.1 | U | | < EQL | | | | | ug/L |
| | alpha-Chlordane | <.1 | U | | < EQL | | | | | ug/L |
| | beta-Benzene hexachloride | <.1 | U | | < EQL | <.098 | U | | < EQL | ug/L |
| | cis-1,2-Dichloroethylene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | cis-1,3-Dichloropropene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | delta-Benzene hexachloride | <.1 | U | | < EQL | | | | | ug/L |
| | gamma-Chlordane | <.1 | U | | < EQL | | | | | ug/L |
| | m-Cresol (3-Methylphenol) | | | | | | | | | |
| | m-Nitroaniline | <26 | U | | < EQL | | | | | ug/L |
| | o-Cresol (2-Methylphenol) | <11 | U | | < EQL | | | | | ug/L |
| | o-Nitroaniline | <26 | U | | < EQL | | | | | ug/L |
| | o-Toluidine | <11 | U | | < EQL | | | | | ug/L |
| | p,p'-DDD | <.21 | U | | < EQL | | | | | ug/L |
| | p,p'-DDE | <.21 | U | | < EQL | | | | | ug/L |
| | p,p'-DDT | <.21 | U | | < EQL | | | | | ug/L |
| | p-Cresol (4-Methylphenol) | <10 | U | | < EQL | | | | | ug/L |
| | p-Dimethylaminoazobenzene | <11 | U | | < EQL | | | | | ug/L |
| | p-Nitroaniline | <11 | U | | < EQL | | | | | ug/L |
| | p-Phenylenediamine | <11 | U | | < EQL | | | | | ug/L |
| | pH | 5.72 | J | Q | NDD | 5.11 | J | Q | NDD | pH |
| | trans-1,2-Dichloroethylene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | trans-1,3-Dichloropropene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | trans-1,4-Dichloro-2-butene | <20 | U | | < EQL | <20 | U | | < EQL | ug/L |

Notes: Concentrations in bold italics exceed the groundwater protection standards listed in Appendix A. Bold Italics were not assigned if the result was qualified with a "J", "R", "L", or "U" qualifier.

+ = exceeded the Groundwater Protection Standards listed in Appendix A.

Table B. Assessment Monitoring Results for Groundwater Wells (Cont.)

WELL LFW 43D

| <u>SRS Coord.</u> | <u>Lat/Longitude</u> | <u>Screen Zone Elevation</u> | <u>Top of Casing</u> | <u>Casing</u> | <u>Pump</u> | <u>Screen Zone</u> |
|------------------------|----------------------------------|------------------------------|----------------------|---------------|-------------|--------------------|
| N 86443.2 E 45244.5 | 33.29077 Deg N -81.7156 Deg W | 170.9 - 150.9 ft msl | 202.9 ft msl | 4 " PVC | S | Upper |

SAMPLE DATE 02/28/00 10/02/00

FIELD DATA

| <u>Parameter</u> | <u>1st Half</u> | <u>2nd Half</u> | <u>Unit</u> |
|---------------------------------|-----------------|-----------------|-------------|
| Water Elevation | 163.8 | 162.88 | ft msl |
| pH | 5 | 5.2 | pH |
| Sp. Conductance | 11 | 11 | uS/cm |
| Water temperature | 17.8 | 18.8 | deg. C |
| Alkalinity as CaCO ₃ | 3 | 0 | mg/L |
| Turbidity | 1.2 | 1.3 | NTU |
| Volumes purged | 6.12 | 4.01 | well volume |
| Sampling code | | | |

ANALYTICAL DATA

Groundwater Protection Standard

| <u>ST</u> | <u>Parameter</u> | <u>1st Half</u> | <u>CLP</u> | <u>EPA</u> | <u>Filt.</u> | <u>2nd Half</u> | <u>CLP</u> | <u>EPA</u> | <u>Filt.</u> | <u>Unit</u> |
|-----------|--------------------------------|-----------------|------------|------------|--------------|-----------------|------------|------------|--------------|-------------|
| | 1,1,1,2-Tetrachloroethane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,1,1-Trichloroethane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,1,2,2-Tetrachloroethane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,1,2-Trichloroethane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,1-Dichloroethane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,1-Dichloroethylene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,1-Dichloropropene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,2,3-Trichloropropane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,2,4,5-Tetrachlorobenzene | <52 | U | | < EQL | | | | | ug/L |
| | 1,2,4-Trichlorobenzene | <11 | U | | < EQL | | | | | ug/L |
| | 1,2-Dibromo-3-chloropropane | <10 | U | | < EQL | <10 | U | | < EQL | ug/L |
| | 1,2-Dibromoethane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,2-Dichlorobenzene | <11 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,2-Dichloroethane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,2-Dichloropropane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,3,5-Trinitrobenzene | <10 | U | | < EQL | | | | | ug/L |
| | 1,3-Dichlorobenzene | <11 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,3-Dichloropropane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,3-Dinitrobenzene | <50 | U | | < EQL | | | | | ug/L |
| | 1,4-Dichlorobenzene | <11 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,4-Naphthoquinone | <11 | U | | < EQL | | | | | ug/L |
| | 1-Naphthylamine | <11 | U | | < EQL | | | | | ug/L |
| | 2,2-Dichloropropane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 2,2-Oxybis(1-chloropropane) | <11 | U | | < EQL | | | | | ug/L |
| | 2,3,4,6-Tetrachlorophenol | <11 | U | | < EQL | | | | | ug/L |
| | 2,4,5-T | | | | | | | | | |
| | 2,4,5-TP (Silvex) | | | | | | | | | |
| | 2,4,5-Trichlorophenol | <11 | U | | < EQL | | | | | ug/L |
| | 2,4,6-Trichlorophenol | <26 | U | | < EQL | | | | | ug/L |
| | 2,4-Dichlorophenol | <11 | U | | < EQL | | | | | ug/L |
| | 2,4-Dichlorophenoxyacetic acid | | | | | | | | | |
| | 2,4-Dimethyl phenol | <11 | U | | < EQL | | | | | ug/L |
| | 2,4-Dinitrophenol | <26 | U | | < EQL | | | | | ug/L |
| | 2,4-Dinitrotoluene | <11 | U | | < EQL | | | | | ug/L |
| | 2,6-Dichlorophenol | <25 | U | | < EQL | | | | | ug/L |
| | 2,6-Dinitrotoluene | <11 | U | | < EQL | | | | | ug/L |
| | 2-Acetylaminofluorene | <11 | U | | < EQL | | | | | ug/L |
| | 2-Chloronaphthalene | <11 | U | | < EQL | | | | | ug/L |
| | 2-Chlorophenol | <11 | U | | < EQL | | | | | ug/L |
| | 2-Hexanone | <20 | U | | < EQL | <20 | U | | < EQL | ug/L |
| | 2-Methyl-4,6-dinitrophenol | <26 | U | | < EQL | | | | | ug/L |
| | 2-Methylnaphthalene | <11 | U | | < EQL | | | | | ug/L |
| | 2-Naphthylamine | <11 | U | | < EQL | | | | | ug/L |
| | 2-Nitrophenol | <11 | U | | < EQL | | | | | ug/L |
| | 2-sec-Butyl-4,6-dinitrophenol | | | | | | | | | |
| | 3,3"-Dichlorobenzidine | <11 | U | | < EQL | | | | | ug/L |
| | 3,3"-Dimethylbenzidine | <21 | U | | < EQL | | | | | ug/L |
| | 3-Methylcholanthrene | <11 | U | | < EQL | | | | | ug/L |
| | 4-Aminobiphenyl | <11 | U | | < EQL | | | | | ug/L |
| | 4-Bromophenyl phenyl ether | <11 | U | | < EQL | | | | | ug/L |
| | 4-Chloro-m-cresol | <11 | U | | < EQL | | | | | ug/L |

Notes: Concentrations in bold italics exceed the groundwater protection standards listed in Appendix A. Bold italics were not assigned if the result was qualified with a "J", "R", "L", or "U" qualifier.

+ = exceeded the Groundwater Protection Standards listed in Appendix A.

Table B. Assessment Monitoring Results for Groundwater Wells (Cont.)

WELL LFW 43D

ANALYTICAL DATA

Groundwater Protection Standard

| ST | Parameter | 1st Half | CLP | EPA | Filt. | 2nd Half | CLP | EPA | Filt. | Unit |
|----|--------------------------------------|----------|-----|-----|-------|----------|-----|-----|-------|------|
| | 4-Chloroaniline | <11 | U | | < EQL | | | | | ug/L |
| | 4-Chlorophenyl phenyl ether | <11 | U | | < EQL | | | | | ug/L |
| | 4-Nitrophenol | <26 | U | | < EQL | | | | | ug/L |
| | 5-Nitro-o-toluidine | <11 | U | | < EQL | | | | | ug/L |
| | 7,12-Dimethylbenz(a)anthracene | <11 | U | | < EQL | | | | | ug/L |
| | Acenaphthene | <11 | U | | < EQL | | | | | ug/L |
| | Acenaphthylene | <11 | U | | < EQL | | | | | ug/L |
| | Acetone | <20 | U | | < EQL | <20 | U | | < EQL | ug/L |
| | Acetonitrile (Methyl cyanide) | <200 | U | | < EQL | <200 | U | | < EQL | ug/L |
| | Acetophenone | <11 | U | | < EQL | | | | | ug/L |
| | Acrolein | <50 | U | | < EQL | <50 | U | | < EQL | ug/L |
| | Acrylonitrile | <10 | U | | < EQL | <10 | U | | < EQL | ug/L |
| | Aldrin | <.1 | U | | < EQL | | | | | ug/L |
| | Allyl chloride | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Anthracene | <11 | U | | < EQL | | | | | ug/L |
| | Antimony, total recoverable | <100 | U | | < EQL | <100 | U | | < EQL | ug/L |
| | Arsenic, total recoverable | <10 | U | | < EQL | 5.13 | J | I | NDD | ug/L |
| | Barium, total recoverable | 6.65 | J | I | NDD | 6.2 | J | I | NDD | ug/L |
| | Benzene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Benzo(a)anthracene | <11 | U | | < EQL | | | | | ug/L |
| | Benzo(a)pyrene | <11 | U | | < EQL | | | | | ug/L |
| | Benzo(b)fluoranthene | <11 | U | | < EQL | | | | | ug/L |
| | Benzo(g,h,i)perylene | <11 | U | | < EQL | | | | | ug/L |
| | Benzo(k)fluoranthene | <11 | U | | < EQL | | | | | ug/L |
| | Benzyl alcohol | | | | | | | | | |
| | Beryllium, total recoverable | <1 | U | | < EQL | <.12 | JU | | < EQL | ug/L |
| | Bis(2-chloroethoxy) methane | <11 | U | | < EQL | | | | | ug/L |
| | Bis(2-chloroethyl) ether | <11 | U | | < EQL | | | | | ug/L |
| | Bis(2-ethylhexyl) phthalate | <11 | U | | < EQL | | | | | ug/L |
| | Bromochloromethane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Bromodichloromethane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Bromoform | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Bromomethane (Methyl bromide) | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Butylbenzyl phthalate | <11 | U | | < EQL | | | | | ug/L |
| | Cadmium, total recoverable | <10 | U | | < EQL | <10 | U | | < EQL | ug/L |
| | Carbon disulfide | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Carbon tetrachloride | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Chlorobenzene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Chlorobenzilate | <11 | U | | < EQL | | | | | ug/L |
| | Chloroethane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Chloroethene (Vinyl chloride) | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Chloroform | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Chloromethane (Methyl chloride) | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Chloroprene | <20 | U | | < EQL | <20 | U | | < EQL | ug/L |
| | Chromium, total recoverable | <3.55 | JU | V | < EQL | <10 | U | | < EQL | ug/L |
| | Chrysene | <11 | U | | < EQL | | | | | ug/L |
| | Cobalt, total recoverable | <20 | U | | < EQL | <20 | U | | < EQL | ug/L |
| | Copper, total recoverable | <11.9 | U | V | < EQL | 7.62 | J | I | NDD | ug/L |
| | Cyanide | | | | | | | | | |
| | Di-n-butyl phthalate | <11 | U | | < EQL | <9.5 | U | | < EQL | ug/L |
| | Di-n-octyl phthalate | <11 | U | | < EQL | | | | | ug/L |
| | Diallate | <11 | U | | < EQL | | | | | ug/L |
| | Dibenz(a,h)anthracene | <11 | U | | < EQL | | | | | ug/L |
| | Dibenzofuran | <11 | U | | < EQL | | | | | ug/L |
| | Dibromochloromethane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Dibromomethane (Methylene bromide) | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Dichlorodifluoromethane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Dichloromethane (Methylene chloride) | <4.9 | U | V | < EQL | <10 | U | | < EQL | ug/L |
| | Dieldrin | <.21 | U | | < EQL | | | | | ug/L |
| | Diethyl phthalate | <11 | U | | < EQL | | | | | ug/L |
| | Dimethoate | <11 | U | | < EQL | | | | | ug/L |
| | Dimethyl phthalate | <11 | U | | < EQL | | | | | ug/L |
| | Diphenylamine | <11 | U | | < EQL | | | | | ug/L |
| | Disulfoton | <11 | U | | < EQL | | | | | ug/L |
| | Endosulfan I | <.1 | U | | < EQL | | | | | ug/L |
| | Endosulfan II | <.21 | U | | < EQL | | | | | ug/L |
| | Endosulfan sulfate | <.21 | U | | < EQL | | | | | ug/L |
| | Endrin | <.21 | U | | < EQL | | | | | ug/L |

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+ = exceeded the Groundwater Protection Standards listed in Appendix A.

Table B. Assessment Monitoring Results for Groundwater Wells (Cont.)

WELL LFW 43D

ANALYTICAL DATA

Groundwater Protection Standard

| ST | Parameter | 1st Half | CLP | EPA | Filt. | 2nd Half | CLP | EPA | Filt. | Unit |
|----|---------------------------------|----------|-----|-----|-------|----------|-----|-----|-------|-------|
| | Endrin aldehyde | <.21 | U | | < EQL | | | | | ug/L |
| | Ethyl methacrylate | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Ethyl methanesulfonate | <11 | U | | < EQL | | | | | ug/L |
| | Ethylbenzene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Famphur | <200 | U | | < EQL | | | | | ug/L |
| | Fluoranthene | <11 | U | | < EQL | | | | | ug/L |
| | Fluorene | <11 | U | | < EQL | | | | | ug/L |
| | Heptachlor | <.1 | U | | < EQL | | | | | ug/L |
| | Heptachlor epoxide | <.1 | U | | < EQL | | | | | ug/L |
| | Hexachlorobenzene | <11 | U | | < EQL | | | | | ug/L |
| | Hexachlorobutadiene | <11 | U | | < EQL | | | | | ug/L |
| | Hexachlorocyclopentadiene | <11 | U | | < EQL | | | | | ug/L |
| | Hexachloroethane | <11 | U | | < EQL | | | | | ug/L |
| | Hexachloropropene | <52 | U | | < EQL | | | | | ug/L |
| | Indeno(1,2,3-c,d)pyrene | <11 | U | | < EQL | | | | | ug/L |
| | Iodomethane (Methyl iodide) | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Isobutyl alcohol | <500 | U | | < EQL | <500 | U | | < EQL | ug/L |
| | Isodrin | <11 | U | | < EQL | | | | | ug/L |
| | Isophorone | <11 | U | | < EQL | | | | | ug/L |
| | Isosafrole | <11 | U | | < EQL | | | | | ug/L |
| | Kepone | <11 | U | | < EQL | | | | | ug/L |
| | Lead, total recoverable | <10 | U | | < EQL | <10 | U | | < EQL | ug/L |
| | Lindane | <.1 | U | | < EQL | | | | | ug/L |
| | Mercury, total recoverable | <.5 | U | | < EQL | <.5 | U | | < EQL | ug/L |
| | Methacrylonitrile | <200 | U | | < EQL | <200 | U | | < EQL | ug/L |
| | Methapyrilene | <11 | U | | < EQL | | | | | ug/L |
| | Methoxychlor | <.1 | U | | < EQL | | | | | ug/L |
| | Methyl ethyl ketone | <20 | U | | < EQL | <20 | U | | < EQL | ug/L |
| | Methyl isobutyl ketone | <10 | U | | < EQL | <10 | U | | < EQL | ug/L |
| | Methyl methacrylate | <20 | U | | < EQL | <20 | U | | < EQL | ug/L |
| | Methyl methanesulfonate | <11 | U | | < EQL | | | | | ug/L |
| | N-Nitrosodi-n-butylamine | <11 | U | | < EQL | | | | | ug/L |
| | N-Nitrosodiethylamine | <11 | U | | < EQL | | | | | ug/L |
| | N-Nitrosodimethylamine | <26 | U | | < EQL | | | | | ug/L |
| | N-Nitrosodiphenylamine | <11 | U | | < EQL | | | | | ug/L |
| | N-Nitrosodipropylamine | <11 | U | | < EQL | | | | | ug/L |
| | N-Nitrosomethylethylamine | <11 | U | | < EQL | | | | | ug/L |
| | N-Nitrosopiperidine | <11 | U | | < EQL | | | | | ug/L |
| | N-Nitrosopyrrolidine | <11 | U | | < EQL | | | | | ug/L |
| | Naphthalene | <11 | U | | < EQL | | | | | ug/L |
| | Nickel, total recoverable | <50 | U | | < EQL | <50 | U | | < EQL | ug/L |
| | Nitrobenzene | <11 | U | | < EQL | | | | | ug/L |
| | O,O,O-Triethyl phosphorothioate | <11 | U | | < EQL | | | | | ug/L |
| | PCB | | | | | | | | | |
| | PCB 1016 | | | | | | | | | |
| | PCB 1232 | | | | | | | | | |
| | PCB 1242 | | | | | | | | | |
| | PCB 1248 | | | | | | | | | |
| | PCB 1254 | | | | | | | | | |
| | PCB 1260 | | | | | | | | | |
| | Parathion | <11 | U | | < EQL | | | | | ug/L |
| | Parathion methyl | <11 | U | | < EQL | | | | | ug/L |
| | Pentachlorobenzene | <52 | U | | < EQL | | | | | ug/L |
| | Pentachloronitrobenzene | <11 | U | | < EQL | | | | | ug/L |
| | Pentachlorophenol | <26 | U | | < EQL | | | | | ug/L |
| | Phenacetin | <11 | U | | < EQL | | | | | ug/L |
| | Phenanthrene | <11 | U | | < EQL | | | | | ug/L |
| | Phenol | <11 | U | | < EQL | | | | | ug/L |
| | Phorate | <11 | U | | < EQL | | | | | ug/L |
| | Pronamid | <11 | U | | < EQL | | | | | ug/L |
| | Propionitrile | <200 | U | | < EQL | <200 | U | | < EQL | ug/L |
| | Pyrene | <11 | U | | < EQL | | | | | ug/L |
| | Safrole | <11 | U | | < EQL | | | | | ug/L |
| | Selenium, total recoverable | <10 | U | | < EQL | <10 | U | | < EQL | ug/L |
| | Silver, total recoverable | <7.07 | JU | | < EQL | 3.92 | J | I | NDD | ug/L |
| | Specific conductance | 9.8 | | | < 37 | 8.98 | | | < 37 | uS/cm |
| | Styrene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Sulfide | | | | | | | | | |

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+ = exceeded the Groundwater Protection Standards listed in Appendix A.

Table B. Assessment Monitoring Results for Groundwater Wells (Cont.)

WELL LFW 43D

ANALYTICAL DATA

Groundwater Protection Standard

| ST | Parameter | 1st Half | CLP | EPA | Filt. | 2nd Half | CLP | EPA | Filt. | Unit |
|----|-----------------------------|----------|-----|-----|-------|-------------|-----|-----|-------|------|
| | Tetrachloroethylene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| + | Thallium, total recoverable | <5.98 | JU | | < EQL | 10.1 | | | > 2 | ug/L |
| | Thionazin | <11 | U | | < EQL | | | | | ug/L |
| | Tin, total recoverable | <200 | U | | < EQL | | | | | ug/L |
| | Toluene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Toxaphene | <2.1 | U | | < EQL | | | | | ug/L |
| | Trichloroethylene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Trichlorofluoromethane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Vanadium, total recoverable | <10 | U | | < EQL | <10 | U | | < EQL | ug/L |
| | Vinyl acetate | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Xylenes | <10 | U | | < EQL | <10 | U | | < EQL | ug/L |
| | Zinc, total recoverable | 13.5 | J | I | NDD | <20 | U | | < EQL | ug/L |
| | alpha-Benzene hexachloride | <.1 | U | | < EQL | | | | | ug/L |
| | alpha-Chlordane | <.1 | U | | < EQL | | | | | ug/L |
| | beta-Benzene hexachloride | <.1 | U | | < EQL | <.095 | U | | < EQL | ug/L |
| | cis-1,2-Dichloroethylene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | cis-1,3-Dichloropropene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | delta-Benzene hexachloride | <.1 | U | | < EQL | | | | | ug/L |
| | gamma-Chlordane | <.1 | U | | < EQL | | | | | ug/L |
| | m-Cresol (3-Methylphenol) | | | | | | | | | |
| | m-Nitroaniline | <26 | U | | < EQL | | | | | ug/L |
| | o-Cresol (2-Methylphenol) | <11 | U | | < EQL | | | | | ug/L |
| | o-Nitroaniline | <26 | U | | < EQL | | | | | ug/L |
| | o-Toluidine | <11 | U | | < EQL | | | | | ug/L |
| | p,p"-DDD | <.21 | U | | < EQL | | | | | ug/L |
| | p,p"-DDE | <.21 | U | | < EQL | | | | | ug/L |
| | p,p"-DDT | <.21 | U | | < EQL | | | | | ug/L |
| | p-Cresol (4-Methylphenol) | <10 | U | | < EQL | | | | | ug/L |
| | p-Dimethylaminoazobenzene | <11 | U | | < EQL | | | | | ug/L |
| | p-Nitroaniline | <11 | U | | < EQL | | | | | ug/L |
| | p-Phenylenediamine | <11 | U | | < EQL | | | | | ug/L |
| | pH | 5.75 | J | Q | NDD | 5.13 | J | Q | NDD | pH |
| | trans-1,2-Dichloroethylene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | trans-1,3-Dichloropropene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | trans-1,4-Dichloro-2-butene | <20 | U | | < EQL | <20 | U | | < EQL | ug/L |

Notes: Concentrations in bold italics exceed the groundwater protection standards listed in Appendix A. Bold italics were not assigned if the result was qualified with a "J", "R", "L", or "U" qualifier.

+ = exceeded the Groundwater Protection Standards listed in Appendix A.

Table B. Assessment Monitoring Results for Groundwater Wells (Cont.)

WELL LFW 74C

| SRS Coord. | Lat/Longitude | Screen Zone Elevation | Top of Casing | Casing | Pump | Screen Zone |
|------------------------|----------------------------------|-----------------------|---------------|---------|------|-------------|
| N 85813.8 E 45097.8 | 33.28914 Deg N -81.7147 Deg W | 116 - 101 ft msl | 213.6 ft msl | 4 " PVC | S | Middle |

SAMPLE DATE

03/08/00

10/02/00

FIELD DATA

| Parameter | 1st Half | 2nd Half | Unit |
|---------------------|----------|----------|-------------|
| Water Elevation | 160.8 | 202.18 | ft msl |
| pH | 5 | 5.1 | pH |
| Sp. Conductance | 35 | 35 | uS/cm |
| Water temperature | 19.2 | 21 | deg. C |
| Alkalinity as CaCO3 | 4 | 0 | mg/L |
| Turbidity | .8 | .5 | NTU |
| Volumes purged | 3.89 | 2.57 | well volume |
| Sampling code | | | |

ANALYTICAL DATA

Groundwater Protection Standard

| ST | Parameter | 1st Half | CLP | EPA | Filt. | 2nd Half | CLP | EPA | Filt. | Unit |
|----|--------------------------------|----------|-----|-----|-------|----------|-----|-----|-------|------|
| | 1,1,1,2-Tetrachloroethane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,1,1-Trichloroethane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,1,2,2-Tetrachloroethane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,1,2-Trichloroethane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,1-Dichloroethane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,1-Dichloroethylene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,1-Dichloropropene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,2,3-Trichloropropane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,2,4,5-Tetrachlorobenzene | <53 | JU | Q | < EQL | | | | | ug/L |
| | 1,2,4-Trichlorobenzene | <11 | JU | Q | < EQL | | | | | ug/L |
| | 1,2-Dibromo-3-chloropropane | <10 | U | | < EQL | <10 | U | | < EQL | ug/L |
| | 1,2-Dibromoethane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,2-Dichlorobenzene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,2-Dichloroethane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,2-Dichloropropane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,3,5-Trinitrobenzene | <10 | JU | Q | < EQL | | | | | ug/L |
| | 1,3-Dichlorobenzene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,3-Dichloropropane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,3-Dinitrobenzene | <50 | JU | Q | < EQL | | | | | ug/L |
| | 1,4-Dichlorobenzene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,4-Naphthoquinone | <11 | JU | Q | < EQL | | | | | ug/L |
| | 1-Naphthylamine | <11 | JU | Q | < EQL | | | | | ug/L |
| | 2,2-Dichloropropane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 2,2-Oxybis(1-chloropropane) | <11 | JU | Q | < EQL | | | | | ug/L |
| | 2,3,4,6-Tetrachlorophenol | <11 | JU | Q | < EQL | | | | | ug/L |
| | 2,4,5-T | <2 | U | Q | < EQL | | | | | ug/L |
| | 2,4,5-TP (Silvex) | <2 | U | Q | < EQL | | | | | ug/L |
| | 2,4,5-Trichlorophenol | <11 | JU | Q | < EQL | | | | | ug/L |
| | 2,4,6-Trichlorophenol | <26 | JU | Q | < EQL | | | | | ug/L |
| | 2,4-Dichlorophenol | <11 | JU | Q | < EQL | | | | | ug/L |
| | 2,4-Dichlorophenoxyacetic acid | <2 | U | Q | < EQL | | | | | ug/L |
| | 2,4-Dimethyl phenol | <11 | JU | Q | < EQL | | | | | ug/L |
| | 2,4-Dinitrophenol | <26 | JU | Q | < EQL | | | | | ug/L |
| | 2,4-Dinitrotoluene | <11 | JU | Q | < EQL | | | | | ug/L |
| | 2,6-Dichlorophenol | <25 | JU | Q | < EQL | | | | | ug/L |
| | 2,6-Dinitrotoluene | <11 | JU | Q | < EQL | | | | | ug/L |
| | 2-Acetylaminofluorene | <11 | JU | Q | < EQL | | | | | ug/L |
| | 2-Chloronaphthalene | <11 | JU | Q | < EQL | | | | | ug/L |
| | 2-Chlorophenol | <11 | JU | Q | < EQL | | | | | ug/L |
| | 2-Hexanone | <20 | U | | < EQL | <20 | U | | < EQL | ug/L |
| | 2-Methyl-4,6-dinitrophenol | <26 | JU | Q | < EQL | | | | | ug/L |
| | 2-Methylnaphthalene | <11 | JU | Q | < EQL | | | | | ug/L |
| | 2-Naphthylamine | <11 | JU | Q | < EQL | | | | | ug/L |
| | 2-Nitrophenol | <11 | JU | Q | < EQL | | | | | ug/L |
| | 2-sec-Butyl-4,6-dinitrophenol | <2 | U | Q | < EQL | | | | | ug/L |
| | 3,3'-Dichlorobenzidine | <11 | JU | Q | < EQL | | | | | ug/L |
| | 3,3'-Dimethylbenzidine | <21 | JU | Q | < EQL | | | | | ug/L |
| | 3-Methylcholanthrene | <11 | JU | Q | < EQL | | | | | ug/L |
| | 4-Aminobiphenyl | <11 | JU | Q | < EQL | | | | | ug/L |
| | 4-Bromophenyl phenyl ether | <11 | JU | Q | < EQL | | | | | ug/L |
| | 4-Chloro-m-cresol | <11 | JU | Q | < EQL | | | | | ug/L |

Notes: Concentrations in bold italics exceed the groundwater protection standards listed in Appendix A. Bold italics were not assigned if the result was qualified with a "J", "R", "L", or "U" qualifier.

+ = exceeded the Groundwater Protection Standards listed in Appendix A.

Table B. Assessment Monitoring Results for Groundwater Wells (Cont.)

WELL LFW 74C

ANALYTICAL DATA

Groundwater Protection Standard

| ST | Parameter | 1st Half | CLP | EPA | Filt. | 2nd Half | CLP | EPA | Filt. | Unit |
|----|--------------------------------------|----------|-----|-----|-------|----------|-----|-----|-------|------|
| | 4-Chloroaniline | <11 | JU | Q | < EQL | | | | | ug/L |
| | 4-Chlorophenyl phenyl ether | <11 | JU | Q | < EQL | | | | | ug/L |
| | 4-Nitrophenol | <26 | JU | Q | < EQL | | | | | ug/L |
| | 5-Nitro-o-toluidine | <11 | JU | Q | < EQL | | | | | ug/L |
| | 7,12-Dimethylbenz(a)anthracene | <11 | JU | Q | < EQL | | | | | ug/L |
| | Acenaphthene | <11 | JU | Q | < EQL | | | | | ug/L |
| | Acenaphthylene | <11 | JU | Q | < EQL | | | | | ug/L |
| | Acetone | <20 | U | | < EQL | <20 | U | | < EQL | ug/L |
| | Acetonitrile (Methyl cyanide) | <200 | U | | < EQL | <200 | U | | < EQL | ug/L |
| | Acetophenone | <11 | JU | Q | < EQL | | | | | ug/L |
| | Acrolein | <50 | U | | < EQL | <50 | U | | < EQL | ug/L |
| | Acrylonitrile | <10 | U | | < EQL | <10 | U | | < EQL | ug/L |
| | Aldrin | <.1 | U | | < EQL | | | | | ug/L |
| | Allyl chloride | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Anthracene | <11 | JU | Q | < EQL | | | | | ug/L |
| | Antimony, total recoverable | <100 | U | | < EQL | <100 | U | | < EQL | ug/L |
| | Arsenic, total recoverable | <10 | U | | < EQL | <10 | U | | < EQL | ug/L |
| | Barium, total recoverable | 5.97 | J | I | NDD | 5.65 | J | I | NDD | ug/L |
| | Benzene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Benzo(a)anthracene | <11 | JU | Q | < EQL | | | | | ug/L |
| | Benzo(a)pyrene | <11 | JU | Q | < EQL | | | | | ug/L |
| | Benzo(b)fluoranthene | <11 | JU | Q | < EQL | | | | | ug/L |
| | Benzo(g,h,i)perylene | <11 | JU | Q | < EQL | | | | | ug/L |
| | Benzo(k)fluoranthene | <11 | JU | Q | < EQL | | | | | ug/L |
| | Benzyl alcohol | | | | | | | | | |
| | Beryllium, total recoverable | <1 | U | | < EQL | <.13 | JU | | < EQL | ug/L |
| | Bis(2-chloroethoxy) methane | <11 | JU | Q | < EQL | | | | | ug/L |
| | Bis(2-chloroethyl) ether | <11 | JU | Q | < EQL | | | | | ug/L |
| | Bis(2-ethylhexyl) phthalate | <11 | JU | Q | < EQL | | | | | ug/L |
| | Bromochloromethane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Bromodichloromethane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Bromoform | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Bromomethane (Methyl bromide) | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Butylbenzyl phthalate | <11 | JU | Q | < EQL | | | | | ug/L |
| | Cadmium, total recoverable | <10 | U | | < EQL | <10 | U | | < EQL | ug/L |
| | Carbon disulfide | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Carbon tetrachloride | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Chlorobenzene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Chlorobenzilate | <11 | JU | Q | < EQL | | | | | ug/L |
| | Chloroethane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Chloroethene (Vinyl chloride) | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Chloroform | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Chloromethane (Methyl chloride) | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Chloroprene | <20 | U | | < EQL | <20 | U | | < EQL | ug/L |
| | Chromium, total recoverable | <10 | U | | < EQL | <10 | U | | < EQL | ug/L |
| | Chrysene | <11 | JU | Q | < EQL | | | | | ug/L |
| | Cobalt, total recoverable | <20 | U | | < EQL | <20 | U | | < EQL | ug/L |
| | Copper, total recoverable | <20 | U | | < EQL | <2.67 | JU | | < EQL | ug/L |
| | Cyanide | 10 | J | Q | NDD | | | | | ug/L |
| | Di-n-butyl phthalate | <11 | JU | Q | < EQL | <9.6 | U | | < EQL | ug/L |
| | Di-n-octyl phthalate | <11 | JU | Q | < EQL | | | | | ug/L |
| | Diallate | <11 | JU | Q | < EQL | | | | | ug/L |
| | Dibenz(a,h)anthracene | <11 | JU | Q | < EQL | | | | | ug/L |
| | Dibenzofuran | <11 | JU | Q | < EQL | | | | | ug/L |
| | Dibromochloromethane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Dibromomethane (Methylene bromide) | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Dichlorodifluoromethane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Dichloromethane (Methylene chloride) | 1.5 | J | I | NDD | <10 | U | | < EQL | ug/L |
| | Dieldrin | <.21 | U | | < EQL | | | | | ug/L |
| | Diethyl phthalate | <11 | JU | Q | < EQL | | | | | ug/L |
| | Dimethoate | <11 | JU | Q | < EQL | | | | | ug/L |
| | Dimethyl phthalate | <11 | JU | Q | < EQL | | | | | ug/L |
| | Diphenylamine | <11 | JU | Q | < EQL | | | | | ug/L |
| | Disulfoton | <11 | JU | Q | < EQL | | | | | ug/L |
| | Endosulfan I | <.1 | U | | < EQL | | | | | ug/L |
| | Endosulfan II | <.21 | U | | < EQL | | | | | ug/L |
| | Endosulfan sulfate | <.21 | U | | < EQL | | | | | ug/L |
| | Endrin | <.21 | U | | < EQL | | | | | ug/L |

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+ = exceeded the Groundwater Protection Standards listed in Appendix A.

Table B. Assessment Monitoring Results for Groundwater Wells (Cont.)

WELL LFW 74C

ANALYTICAL DATA

Groundwater Protection Standard

| ST | Parameter | 1st Half | CLP | EPA | Filt. | 2nd Half | CLP | EPA | Filt. | Unit |
|----|---------------------------------|----------|-----|-----|-------|----------|-----|-----|-------|-------|
| | Endrin aldehyde | <.21 | U | | < EQL | | | | | ug/L |
| | Ethyl methacrylate | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Ethyl methanesulfonate | <11 | JU | Q | < EQL | | | | | ug/L |
| | Ethylbenzene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Famphur | <200 | JU | Q | < EQL | | | | | ug/L |
| | Fluoranthene | <11 | JU | Q | < EQL | | | | | ug/L |
| | Fluorene | <11 | JU | Q | < EQL | | | | | ug/L |
| | Heptachlor | <.1 | U | | < EQL | | | | | ug/L |
| | Heptachlor epoxide * | <.1 | U | | < EQL | | | | | ug/L |
| | Hexachlorobenzene | <11 | JU | Q | < EQL | | | | | ug/L |
| | Hexachlorobutadiene | <11 | JU | Q | < EQL | | | | | ug/L |
| | Hexachlorocyclopentadiene | <11 | JU | Q | < EQL | | | | | ug/L |
| | Hexachloroethane | <11 | JU | Q | < EQL | | | | | ug/L |
| | Hexachloropropene | <53 | JU | Q | < EQL | | | | | ug/L |
| | Indeno(1,2,3-c,d)pyrene | <11 | JU | Q | < EQL | | | | | ug/L |
| | Iodomethane (Methyl iodide) | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Isobutyl alcohol | <500 | U | | < EQL | <500 | U | | < EQL | ug/L |
| | Isodrin | <11 | JU | Q | < EQL | | | | | ug/L |
| | Isophorone | <11 | JU | Q | < EQL | | | | | ug/L |
| | Isosafrole | <11 | JU | Q | < EQL | | | | | ug/L |
| | Kepone | <11 | JU | Q | < EQL | | | | | ug/L |
| | Lead, total recoverable | 4.21 | J | I | NDD | <1.45 | JU | | < EQL | ug/L |
| | Lindane | <.1 | U | | < EQL | | | | | ug/L |
| | Mercury, total recoverable | <.154 | U | V | < EQL | <.5 | U | | < EQL | ug/L |
| | Methacrylonitrile | <200 | U | | < EQL | <200 | U | | < EQL | ug/L |
| | Methapyriene | <11 | JU | Q | < EQL | | | | | ug/L |
| | Methoxychlor | <.1 | U | | < EQL | | | | | ug/L |
| | Methyl ethyl ketone | <20 | U | | < EQL | <20 | U | | < EQL | ug/L |
| | Methyl isobutyl ketone | <10 | U | | < EQL | <10 | U | | < EQL | ug/L |
| | Methyl methacrylate | <20 | U | | < EQL | <20 | U | | < EQL | ug/L |
| | Methyl methanesulfonate | <11 | JU | Q | < EQL | | | | | ug/L |
| | N-Nitrosodi-n-butylamine | <11 | JU | Q | < EQL | | | | | ug/L |
| | N-Nitrosodiethylamine | <11 | JU | Q | < EQL | | | | | ug/L |
| | N-Nitrosodimethylamine | <26 | JU | Q | < EQL | | | | | ug/L |
| | N-Nitrosodiphenylamine | <11 | JU | Q | < EQL | | | | | ug/L |
| | N-Nitrosodipropylamine | <11 | JU | Q | < EQL | | | | | ug/L |
| | N-Nitrosomethylethylamine | <11 | JU | Q | < EQL | | | | | ug/L |
| | N-Nitrosopiperidine | <11 | JU | Q | < EQL | | | | | ug/L |
| | N-Nitrosopyrrolidine | <11 | JU | Q | < EQL | | | | | ug/L |
| | Naphthalene | <11 | JU | Q | < EQL | | | | | ug/L |
| | Nickel, total recoverable | <50 | U | | < EQL | <50 | U | | < EQL | ug/L |
| | Nitrobenzene | <11 | JU | Q | < EQL | | | | | ug/L |
| | O,O,O-Triethyl phosphorothioate | <11 | JU | Q | < EQL | | | | | ug/L |
| | PCB | | | | | | | | | |
| | PCB 1016 | <.1 | U | | < EQL | | | | | ug/L |
| | PCB 1232 | <.1 | U | | < EQL | | | | | ug/L |
| | PCB 1242 | <2.1 | U | | < EQL | | | | | ug/L |
| | PCB 1248 | <.1 | U | | < EQL | | | | | ug/L |
| | PCB 1254 | <.1 | U | | < EQL | | | | | ug/L |
| | PCB 1260 | <.1 | U | | < EQL | | | | | ug/L |
| | Parathion | <11 | JU | Q | < EQL | | | | | ug/L |
| | Parathion methyl | <11 | JU | Q | < EQL | | | | | ug/L |
| | Pentachlorobenzene | <53 | JU | Q | < EQL | | | | | ug/L |
| | Pentachloronitrobenzene | <11 | JU | Q | < EQL | | | | | ug/L |
| | Pentachlorophenol | <26 | JU | Q | < EQL | | | | | ug/L |
| | Phenacetin | <11 | JU | Q | < EQL | | | | | ug/L |
| | Phenanthrene | <11 | JU | Q | < EQL | | | | | ug/L |
| | Phenol | <11 | JU | Q | < EQL | | | | | ug/L |
| | Phorate | <11 | JU | Q | < EQL | | | | | ug/L |
| | Pronamid | <11 | JU | Q | < EQL | | | | | ug/L |
| | Propionitrile | <200 | U | | < EQL | <200 | U | | < EQL | ug/L |
| | Pyrene | <11 | JU | Q | < EQL | | | | | ug/L |
| | Safrole | <11 | JU | Q | < EQL | | | | | ug/L |
| | Selenium, total recoverable | <10 | U | | < EQL | <10 | U | | < EQL | ug/L |
| | Silver, total recoverable | <20 | U | | < EQL | <20 | U | | < EQL | ug/L |
| | Specific conductance | 33.4 | | | < 37 | 32.2 | | | < 37 | uS/cm |
| | Styrene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Sulfide | <1000 | JU | Q | < EQL | | | | | ug/L |

Notes: Concentrations in bold italics exceed the groundwater protection standards listed in Appendix A. Bold italics were not assigned if the result was qualified with a "J", "R", "L", or "U" qualifier.

+ = exceeded the Groundwater Protection Standards listed in Appendix A.

Table B. Assessment Monitoring Results for Groundwater Wells (Cont.)

WELL LFW 74C

ANALYTICAL DATA

Groundwater Protection Standard

| ST | Parameter | 1st Half | CLP | EPA | Filt. | 2nd Half | CLP | EPA | Filt. | Unit |
|----|-----------------------------|----------|-----|-----|-------|----------|-----|-----|-------|------|
| | Tetrachloroethylene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Thallium, total recoverable | <10 | U | | < EQL | <10 | U | | < EQL | ug/L |
| | Thionazin | <11 | JU | Q | < EQL | | | | | ug/L |
| | Tin, total recoverable | <200 | U | | < EQL | | | | | ug/L |
| | Toluene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Toxaphene | <2.1 | U | | < EQL | | | | | ug/L |
| | Trichloroethylene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Trichlorofluoromethane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Vanadium, total recoverable | <10 | U | | < EQL | <10 | U | | < EQL | ug/L |
| | Vinyl acetate | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Xylenes | <10 | U | | < EQL | <10 | U | | < EQL | ug/L |
| | Zinc, total recoverable | 14.8 | J | I | NDD | <20 | U | | < EQL | ug/L |
| | alpha-Benzene hexachloride | <.1 | U | | < EQL | | | | | ug/L |
| | alpha-Chlordane | <.1 | U | | < EQL | | | | | ug/L |
| | beta-Benzene hexachloride | <.1 | U | | < EQL | <.1 | U | | < EQL | ug/L |
| | cis-1,2-Dichloroethylene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | cis-1,3-Dichloropropene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | delta-Benzene hexachloride | <.1 | U | | < EQL | | | | | ug/L |
| | gamma-Chlordane | <.1 | U | | < EQL | | | | | ug/L |
| | m-Cresol (3-Methylphenol) | | | | | | | | | |
| | m-Nitroaniline | <26 | JU | Q | < EQL | | | | | ug/L |
| | o-Cresol (2-Methylphenol) | <11 | JU | Q | < EQL | | | | | ug/L |
| | o-Nitroaniline | <26 | JU | Q | < EQL | | | | | ug/L |
| | o-Toluidine | <11 | JU | Q | < EQL | | | | | ug/L |
| | p,p"-DDD | <.21 | U | | < EQL | | | | | ug/L |
| | p,p"-DDE | <.21 | U | | < EQL | | | | | ug/L |
| | p,p"-DDT | <.21 | U | | < EQL | | | | | ug/L |
| | p-Cresol (4-Methylphenol) | <10 | JU | Q | < EQL | | | | | ug/L |
| | p-Dimethylaminoazobenzene | <11 | JU | Q | < EQL | | | | | ug/L |
| | p-Nitroaniline | <11 | JU | Q | < EQL | | | | | ug/L |
| | p-Phenylenediamine | <11 | JU | Q | < EQL | | | | | ug/L |
| | pH | 5.7 | J | Q | NDD | 4.97 | J | Q | NDD | pH |
| | trans-1,2-Dichloroethylene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | trans-1,3-Dichloropropene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | trans-1,4-Dichloro-2-butene | <20 | U | | < EQL | <20 | U | | < EQL | ug/L |

Notes: Concentrations in bold italics exceed the groundwater protection standards listed in Appendix A. Bold italics were not assigned if the result was qualified with a "J", "R", "L", or "U" qualifier.

+ = exceeded the Groundwater Protection Standards listed in Appendix A.

Table B. Assessment Monitoring Results for Groundwater Wells (Cont.)

WELL LFW 74D

| <u>SRS Coord.</u> | <u>Lat/Longitude</u> | <u>Screen Zone Elevation</u> | <u>Top of Casing</u> | <u>Casing</u> | <u>Pump</u> | <u>Screen Zone</u> |
|----------------------|----------------------------------|------------------------------|----------------------|---------------|-------------|--------------------|
| N 85828.1 E 45098 | 33.28917 Deg N -81.7148 Deg W | 167.7 - 152.7 ft msl | 213.9 ft msl | 4 " PVC | S | Upper |

SAMPLE DATE 03/08/00 10/03/00

FIELD DATA

| <u>Parameter</u> | <u>1st Half</u> | <u>2nd Half</u> | <u>Unit</u> |
|---------------------|-----------------|-----------------|-------------|
| Water Elevation | 160.9 | 160.3 | ft msl |
| pH | 5.2 | 4.9 | pH |
| Sp. Conductance | 25 | 19 | uS/cm |
| Water temperature | 19.9 | 20.3 | deg. C |
| Alkalinity as CaCO3 | 4 | 0 | mg/L |
| Turbidity | 3.2 | 1.8 | NTU |
| Volumes purged | 5.72 | 10.3 | well volume |
| Sampling code | | | |

ANALYTICAL DATA

Groundwater Protection Standard

| <u>SI</u> | <u>Parameter</u> | <u>1st Half</u> | <u>CLP</u> | <u>EPA</u> | <u>Filt.</u> | <u>2nd Half</u> | <u>CLP</u> | <u>EPA</u> | <u>Filt.</u> | <u>Unit</u> |
|-----------|--------------------------------|-----------------|------------|------------|--------------|-----------------|------------|------------|--------------|-------------|
| | 1,1,1,2-Tetrachloroethane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,1,1-Trichloroethane | 3.8 | J | I | NDD | 2.8 | J | IK | NDD | ug/L |
| | 1,1,2,2-Tetrachloroethane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,1,2-Trichloroethane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,1-Dichloroethane | 3 | J | I | NDD | 2.8 | J | IK | NDD | ug/L |
| | 1,1-Dichloroethylene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,1-Dichloropropene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,2,3-Trichloropropane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,2,4,5-Tetrachlorobenzene | <54 | JU | Q | < EQL | | | | | ug/L |
| | 1,2,4-Trichlorobenzene | <11 | JU | Q | < EQL | | | | | ug/L |
| | 1,2-Dibromo-3-chloropropane | <10 | U | | < EQL | <10 | U | | < EQL | ug/L |
| | 1,2-Dibromoethane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,2-Dichlorobenzene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,2-Dichloroethane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,2-Dichloropropane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,3,5-Trinitrobenzene | <10 | JU | Q | < EQL | | | | | ug/L |
| | 1,3-Dichlorobenzene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,3-Dichloropropane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,3-Dinitrobenzene | <50 | JU | Q | < EQL | | | | | ug/L |
| | 1,4-Dichlorobenzene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,4-Naphthoquinone | <11 | JU | Q | < EQL | | | | | ug/L |
| | 1-Naphthylamine | <11 | JU | Q | < EQL | | | | | ug/L |
| | 2,2-Dichloropropane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 2,2-Oxybis(1-chloropropane) | <11 | JU | Q | < EQL | | | | | ug/L |
| | 2,3,4,6-Tetrachlorophenol | <11 | JU | Q | < EQL | | | | | ug/L |
| | 2,4,5-T | <2 | U | Q | < EQL | | | | | ug/L |
| | 2,4,5-TP (Silvex) | <2 | U | Q | < EQL | | | | | ug/L |
| | 2,4,5-Trichlorophenol | <11 | JU | Q | < EQL | | | | | ug/L |
| | 2,4,6-Trichlorophenol | <27 | JU | Q | < EQL | | | | | ug/L |
| | 2,4-Dichlorophenol | <11 | JU | Q | < EQL | | | | | ug/L |
| | 2,4-Dichlorophenoxyacetic acid | <2 | U | Q | < EQL | | | | | ug/L |
| | 2,4-Dimethyl phenol | <11 | JU | Q | < EQL | | | | | ug/L |
| | 2,4-Dinitrophenol | <27 | JU | Q | < EQL | | | | | ug/L |
| | 2,4-Dinitrotoluene | <11 | JU | Q | < EQL | | | | | ug/L |
| | 2,6-Dichlorophenol | <25 | JU | Q | < EQL | | | | | ug/L |
| | 2,6-Dinitrotoluene | <11 | JU | Q | < EQL | | | | | ug/L |
| | 2-Acetylaminofluorene | <11 | JU | Q | < EQL | | | | | ug/L |
| | 2-Chloronaphthalene | <11 | JU | Q | < EQL | | | | | ug/L |
| | 2-Chlorophenol | <11 | JU | Q | < EQL | | | | | ug/L |
| | 2-Hexanone | <20 | U | | < EQL | <20 | U | | < EQL | ug/L |
| | 2-Methyl-4,6-dinitrophenol | <27 | JU | Q | < EQL | | | | | ug/L |
| | 2-Methylnaphthalene | <11 | JU | Q | < EQL | | | | | ug/L |
| | 2-Naphthylamine | <11 | JU | Q | < EQL | | | | | ug/L |
| | 2-Nitrophenol | <11 | JU | Q | < EQL | | | | | ug/L |
| | 2-sec-Butyl-4,6-dinitrophenol | <2 | U | Q | < EQL | | | | | ug/L |
| | 3,3'-Dichlorobenzidine | <11 | JU | Q | < EQL | | | | | ug/L |
| | 3,3'-Dimethylbenzidine | <22 | JU | Q | < EQL | | | | | ug/L |
| | 3-Methylcholanthrene | <11 | JU | Q | < EQL | | | | | ug/L |
| | 4-Aminobiphenyl | <11 | JU | Q | < EQL | | | | | ug/L |
| | 4-Bromophenyl phenyl ether | <11 | JU | Q | < EQL | | | | | ug/L |
| | 4-Chloro-m-cresol | <11 | JU | Q | < EQL | | | | | ug/L |

Notes: Concentrations in bold italics exceed the groundwater protection standards listed in Appendix A. Bold italics were not assigned if the result was qualified with a "J", "R", "L", or "U" qualifier.

+ = exceeded the Groundwater Protection Standards listed in Appendix A.

Table B. Assessment Monitoring Results for Groundwater Wells (Cont.)

WELL LFW 74D

ANALYTICAL DATA

Groundwater Protection Standard

| ST | Parameter | 1st Half | CLP | EPA | Filt. | 2nd Half | CLP | EPA | Filt. | Unit |
|----|--------------------------------------|----------|-----|-----|----------|----------|-----|-----|-------|------|
| | 4-Chloroaniline | <11 | JU | Q | < EQL | | | | | ug/L |
| | 4-Chlorophenyl phenyl ether | <11 | JU | Q | < EQL | | | | | ug/L |
| | 4-Nitrophenol | <27 | JU | Q | < EQL | | | | | ug/L |
| | 5-Nitro-o-toluidine | <11 | JU | Q | < EQL | | | | | ug/L |
| | 7,12-Dimethylbenz(a)anthracene | <11 | JU | Q | < EQL | | | | | ug/L |
| | Acenaphthene | <11 | JU | Q | < EQL | | | | | ug/L |
| | Acenaphthylene | <11 | JU | Q | < EQL | | | | | ug/L |
| | Acetone | 15 | J | I | NDD | 89 | J | K | NDD | ug/L |
| | Acetonitrile (Methyl cyanide) | <200 | U | | < EQL | <200 | U | | < EQL | ug/L |
| | Acetophenone | <11 | JU | Q | < EQL | | | | | ug/L |
| | Acrolein | <50 | U | | < EQL | <50 | U | | < EQL | ug/L |
| | Acrylonitrile | <10 | U | | < EQL | <10 | U | | < EQL | ug/L |
| | Aldrin | <.1 | U | | < EQL | | | | | ug/L |
| | Allyl chloride | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Anthracene | <11 | JU | Q | < EQL | | | | | ug/L |
| | Antimony, total recoverable | <43.2 | JU | | < EQL | <100 | U | | < EQL | ug/L |
| | Arsenic, total recoverable | <10 | U | | < EQL | <10 | U | | < EQL | ug/L |
| | Barium, total recoverable | 7.7 | J | I | NDD | 5.57 | J | I | NDD | ug/L |
| | Benzene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Benzo(a)anthracene | <11 | JU | Q | < EQL | | | | | ug/L |
| | Benzo(a)pyrene | <11 | JU | Q | < EQL | | | | | ug/L |
| | Benzo(b)fluoranthene | <11 | JU | Q | < EQL | | | | | ug/L |
| | Benzo(g,h,i)perylene | <11 | JU | Q | < EQL | | | | | ug/L |
| | Benzo(k)fluoranthene | <11 | JU | Q | < EQL | | | | | ug/L |
| | Benzyl alcohol | | | | | | | | | |
| | Beryllium, total recoverable | <.12 | JU | | < EQL | <.25 | JU | | < EQL | ug/L |
| | Bis(2-chloroethoxy) methane | <11 | JU | Q | < EQL | | | | | ug/L |
| | Bis(2-chloroethyl) ether | <11 | JU | Q | < EQL | | | | | ug/L |
| | Bis(2-ethylhexyl) phthalate | <11 | JU | Q | < EQL | | | | | ug/L |
| | Bromochloromethane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Bromodichloromethane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Bromoform | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Bromomethane (Methyl bromide) | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Butylbenzyl phthalate | <11 | JU | Q | < EQL | | | | | ug/L |
| | Cadmium, total recoverable | <10 | U | | < EQL | <10 | U | | < EQL | ug/L |
| | Carbon disulfide | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Carbon tetrachloride | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Chlorobenzene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Chlorobenzilate | <11 | JU | Q | < EQL | | | | | ug/L |
| | Chloroethane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Chloroethene (Vinyl chloride) | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Chloroform | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Chloromethane (Methyl chloride) | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Chloroprene | <20 | U | | < EQL | <20 | U | | < EQL | ug/L |
| | Chromium, total recoverable | <10 | U | | < EQL | <10 | U | | < EQL | ug/L |
| | Chrysene | <11 | JU | Q | < EQL | | | | | ug/L |
| | Cobalt, total recoverable | <20 | U | | < EQL | <20 | U | | < EQL | ug/L |
| | Copper, total recoverable | <20 | U | | < EQL | <20 | U | | < EQL | ug/L |
| | Cyanide | 10 | R | Q | Rejected | | | | | ug/L |
| | Di-n-butyl phthalate | <11 | JU | Q | < EQL | <9.8 | JU | Q | < EQL | ug/L |
| | Di-n-octyl phthalate | <11 | JU | Q | < EQL | | | | | ug/L |
| | Diallate | <11 | JU | Q | < EQL | | | | | ug/L |
| | Dibenz(a,h)anthracene | <11 | JU | Q | < EQL | | | | | ug/L |
| | Dibenzofuran | <11 | JU | Q | < EQL | | | | | ug/L |
| | Dibromochloromethane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Dibromomethane (Methylene bromide) | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Dichlorodifluoromethane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Dichloromethane (Methylene chloride) | 4.3 | J | I | NDD | <2.8 | U | | < EQL | ug/L |
| | Dieldrin | <.21 | U | | < EQL | | | | | ug/L |
| | Diethyl phthalate | <11 | JU | Q | < EQL | | | | | ug/L |
| | Dimethoate | <11 | JU | Q | < EQL | | | | | ug/L |
| | Dimethyl phthalate | <11 | JU | Q | < EQL | | | | | ug/L |
| | Diphenylamine | <11 | JU | Q | < EQL | | | | | ug/L |
| | Disulfoton | <11 | JU | Q | < EQL | | | | | ug/L |
| | Endosulfan I | <.1 | U | | < EQL | | | | | ug/L |
| | Endosulfan II | <.21 | U | | < EQL | | | | | ug/L |
| | Endosulfan sulfate | <.21 | U | | < EQL | | | | | ug/L |
| | Endrin | <.21 | U | | < EQL | | | | | ug/L |

Notes: Concentrations in bold Italics exceed the groundwater protection standards listed in Appendix A. Bold Italics were not assigned if the result was qualified with a "J", "R", "L", or "U" qualifier.

+ = exceeded the Groundwater Protection Standards listed in Appendix A.

Table B. Assessment Monitoring Results for Groundwater Wells (Cont.)

WELL LFW 74D

ANALYTICAL DATA

Groundwater Protection Standard

| ST | Parameter | 1st Half | CLP | EPA | Filt. | 2nd Half | CLP | EPA | Filt. | Unit |
|----|---------------------------------|----------|-----|-----|----------|----------|-----|-----|-------|-------|
| | Endrin aldehyde | <.21 | U | | < EQL | | | | | ug/L |
| | Ethyl methacrylate | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Ethyl methanesulfonate | <11 | JU | Q | < EQL | | | | | ug/L |
| | Ethylbenzene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Famphur | <200 | JU | Q | < EQL | | | | | ug/L |
| | Fluoranthene | <11 | JU | Q | < EQL | | | | | ug/L |
| | Fluorene | <11 | JU | Q | < EQL | | | | | ug/L |
| | Heptachlor | <.1 | U | | < EQL | | | | | ug/L |
| | Heptachlor epoxide | <.1 | U | | < EQL | | | | | ug/L |
| | Hexachlorobenzene | <11 | JU | Q | < EQL | | | | | ug/L |
| | Hexachlorobutadiene | <11 | JU | Q | < EQL | | | | | ug/L |
| | Hexachlorocyclopentadiene | <11 | JU | Q | < EQL | | | | | ug/L |
| | Hexachloroethane | <11 | JU | Q | < EQL | | | | | ug/L |
| | Hexachloropropene | <54 | JU | Q | < EQL | | | | | ug/L |
| | Indeno(1,2,3-c,d)pyrene | <11 | JU | Q | < EQL | | | | | ug/L |
| | Iodomethane (Methyl iodide) | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Isobutyl alcohol | <500 | U | | < EQL | <500 | U | | < EQL | ug/L |
| | Isodrin | <11 | JU | Q | < EQL | | | | | ug/L |
| | Isophorone | <11 | JU | Q | < EQL | | | | | ug/L |
| | Isosafrole | <11 | JU | Q | < EQL | | | | | ug/L |
| | Kepone | <11 | JU | Q | < EQL | | | | | ug/L |
| | Lead, total recoverable | <10 | U | | < EQL | <1.74 | JU | | < EQL | ug/L |
| | Lindane | <.1 | U | | < EQL | | | | | ug/L |
| | Mercury, total recoverable | .742 | | | < 2 | .494 | J | I | NDD | ug/L |
| | Methacrylonitrile | <200 | U | | < EQL | <200 | U | | < EQL | ug/L |
| | Methapyrene | <11 | JU | Q | < EQL | | | | | ug/L |
| | Methoxychlor | <1 | U | | < EQL | | | | | ug/L |
| | Methyl ethyl ketone | <20 | U | | < EQL | <20 | U | | < EQL | ug/L |
| | Methyl isobutyl ketone | <10 | U | | < EQL | <10 | U | | < EQL | ug/L |
| | Methyl methacrylate | <20 | U | | < EQL | <20 | U | | < EQL | ug/L |
| | Methyl methanesulfonate | <11 | JU | Q | < EQL | | | | | ug/L |
| | N-Nitrosodi-n-butylamine | <11 | JU | Q | < EQL | | | | | ug/L |
| | N-Nitrosodiethylamine | <11 | JU | Q | < EQL | | | | | ug/L |
| | N-Nitrosodimethylamine | <27 | JU | Q | < EQL | | | | | ug/L |
| | N-Nitrosodiphenylamine | <11 | JU | Q | < EQL | | | | | ug/L |
| | N-Nitrosodipropylamine | <11 | JU | Q | < EQL | | | | | ug/L |
| | N-Nitrosomethylethylamine | <11 | JU | Q | < EQL | | | | | ug/L |
| | N-Nitrosopiperidine | <11 | JU | Q | < EQL | | | | | ug/L |
| | N-Nitrosopyrrolidine | <11 | JU | Q | < EQL | | | | | ug/L |
| | Naphthalene | <11 | JU | Q | < EQL | | | | | ug/L |
| | Nickel, total recoverable | <6.34 | JU | | < EQL | <50 | U | | < EQL | ug/L |
| | Nitrobenzene | <11 | JU | Q | < EQL | | | | | ug/L |
| | O,O,O-Triethyl phosphorothioate | <11 | JU | Q | < EQL | | | | | ug/L |
| | PCB | | | | | | | | | |
| | PCB 1016 | <1 | U | | < EQL | | | | | ug/L |
| | PCB 1232 | <1 | U | | < EQL | | | | | ug/L |
| | PCB 1242 | <2.1 | U | | < EQL | | | | | ug/L |
| | PCB 1248 | <1 | U | | < EQL | | | | | ug/L |
| | PCB 1254 | <1 | U | | < EQL | | | | | ug/L |
| | PCB 1260 | <1 | U | | < EQL | | | | | ug/L |
| | Parathion | <11 | JU | Q | < EQL | | | | | ug/L |
| | Parathion methyl | <11 | JU | Q | < EQL | | | | | ug/L |
| | Pentachlorobenzene | <54 | JU | Q | < EQL | | | | | ug/L |
| | Pentachloronitrobenzene | <11 | JU | Q | < EQL | | | | | ug/L |
| | Pentachlorophenol | <27 | JU | Q | < EQL | | | | | ug/L |
| | Phenacetin | <11 | JU | Q | < EQL | | | | | ug/L |
| | Phenanthrene | <11 | JU | Q | < EQL | | | | | ug/L |
| | Phenol | <11 | JU | Q | < EQL | | | | | ug/L |
| | Phorate | <11 | JU | Q | < EQL | | | | | ug/L |
| | Pronamid | <11 | JU | Q | < EQL | | | | | ug/L |
| | Propionitrile | <200 | U | | < EQL | <200 | U | | < EQL | ug/L |
| | Pyrene | <11 | JU | Q | < EQL | | | | | ug/L |
| | Safrole | <11 | JU | Q | < EQL | | | | | ug/L |
| | Selenium, total recoverable | <10 | U | | < EQL | <10 | U | | < EQL | ug/L |
| | Silver, total recoverable | <20 | U | | < EQL | <20 | U | | < EQL | ug/L |
| + | Specific conductance | 20.3 | | | < 37 | 48.5 | | | > 37 | uS/cm |
| | Styrene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Sulfide | 1000 | R | Q | Rejected | | | | | ug/L |

Notes: Concentrations in bold italics exceed the groundwater protection standards listed in Appendix A. Bold italics were not assigned if the result was qualified with a "J", "R", "L", or "U" qualifier.

+ = exceeded the Groundwater Protection Standards listed in Appendix A.

Table B. Assessment Monitoring Results for Groundwater Wells (Cont.)

WELL LFW 74D

ANALYTICAL DATA

Groundwater Protection Standard

| ST | Parameter | 1st Half | CLP | EPA | Filt. | 2nd Half | CLP | EPA | Filt. | Unit |
|----|-----------------------------|-----------|-----|-----|--------|----------|-----|-----|-------|------|
| | Tetrachloroethylene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Thallium, total recoverable | <10 | U | | < EQL | <10 | U | | < EQL | ug/L |
| | Thionazin | <11 | JU | Q | < EQL | | | | | ug/L |
| | Tin, total recoverable | <200 | U | | < EQL | | | | | ug/L |
| | Toluene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Toxaphene | <2.1 | U | | < EQL | | | | | ug/L |
| | Trichloroethylene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Trichlorofluoromethane | 21 | | | > 20 | 16 | J | K | NDD | ug/L |
| | Vanadium, total recoverable | <10 | U | | < EQL | <10 | U | | < EQL | ug/L |
| | Vinyl acetate | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Xylenes | <10 | U | | < EQL | <10 | U | | < EQL | ug/L |
| | Zinc, total recoverable | 23.8 | | | < 29.3 | <4.98 | JU | | < EQL | ug/L |
| | alpha-Benzene hexachloride | <.1 | U | | < EQL | | | | | ug/L |
| | alpha-Chlordane | <.1 | U | | < EQL | | | | | ug/L |
| | beta-Benzene hexachloride | <.1 | U | | < EQL | <.1 | U | | < EQL | ug/L |
| | cis-1,2-Dichloroethylene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | cis-1,3-Dichloropropene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | delta-Benzene hexachloride | <.1 | U | | < EQL | | | | | ug/L |
| | gamma-Chlordane | <.1 | U | | < EQL | | | | | ug/L |
| | m-Cresol (3-Methylphenol) | | | | | | | | | |
| | m-Nitroaniline | <27 | JU | Q | < EQL | | | | | ug/L |
| | o-Cresol (2-Methylphenol) | <11 | JU | Q | < EQL | | | | | ug/L |
| | o-Nitroaniline | <27 | JU | Q | < EQL | | | | | ug/L |
| | o-Toluidine | <11 | JU | Q | < EQL | | | | | ug/L |
| | p,p"-DDD | <.21 | U | | < EQL | | | | | ug/L |
| | p,p"-DDE | <.21 | U | | < EQL | | | | | ug/L |
| | p,p"-DDT | <.21 | U | | < EQL | | | | | ug/L |
| | p-Cresol (4-Methylphenol) | <10 | JU | Q | < EQL | | | | | ug/L |
| | p-Dimethylaminoazobenzene | <11 | JU | Q | < EQL | | | | | ug/L |
| | p-Nitroaniline | <11 | JU | Q | < EQL | | | | | ug/L |
| | p-Phenylenediamine | <11 | JU | Q | < EQL | | | | | ug/L |
| | pH | 5.8 | J | Q | NDD | 4.09 | J | Q | NDD | pH |
| | trans-1,2-Dichloroethylene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | trans-1,3-Dichloropropene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | trans-1,4-Dichloro-2-butene | <20 | U | | < EQL | <20 | U | | < EQL | ug/L |

Notes: Concentrations in bold italics exceed the groundwater protection standards listed in Appendix A. Bold italics were not assigned if the result was qualified with a "J", "R", "L", or "U" qualifier.

+ = exceeded the Groundwater Protection Standards listed in Appendix A.

Table B. Assessment Monitoring Results for Groundwater Wells (Cont.)

WELL LFW 75C

| <u>SRS Coord.</u> | <u>Lat/Longitude</u> | <u>Screen Zone Elevation</u> | <u>Top of Casing</u> | <u>Casing</u> | <u>Pump</u> | <u>Screen Zone</u> |
|----------------------|----------------------------------|------------------------------|----------------------|---------------|-------------|--------------------|
| N 85856.8 E 45357 | 33.28966 Deg N -81.7141 Deg W | 115.6 - 100.6 ft msl | 197.8 ft msl | 4 " PVC | S | Middle |

SAMPLE DATE 03/08/00 10/02/00

FIELD DATA

| <u>Parameter</u> | <u>1st Half</u> | <u>2nd Half</u> | <u>Unit</u> |
|---------------------|-----------------|-----------------|-------------|
| Water Elevation | 160.7 | 160.23 | ft msl |
| pH | 4.9 | 4.9 | pH |
| Sp. Conductance | 31 | 32 | uS/cm |
| Water temperature | 19.6 | 20.6 | deg. C |
| Alkalinity as CaCO3 | 2 | 0 | mg/L |
| Turbidity | .6 | .8 | NTU |
| Volumes purged | 3.87 | 3.95 | well volume |
| Sampling code | | | |

ANALYTICAL DATA

Groundwater Protection Standard

| <u>ST</u> | <u>Parameter</u> | <u>1st Half</u> | <u>CLP</u> | <u>EPA</u> | <u>Filt.</u> | <u>2nd Half</u> | <u>CLP</u> | <u>EPA</u> | <u>Filt.</u> | <u>Unit</u> |
|-----------|--------------------------------|-----------------|------------|------------|--------------|-----------------|------------|------------|--------------|-------------|
| | 1,1,1,2-Tetrachloroethane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,1,1-Trichloroethane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,1,2,2-Tetrachloroethane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,1,2-Trichloroethane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,1-Dichloroethane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,1-Dichloroethylene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,1-Dichloropropene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,2,3-Trichloropropane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,2,4,5-Tetrachlorobenzene | <53 | JU | Q | < EQL | | | | | ug/L |
| | 1,2,4-Trichlorobenzene | <11 | JU | Q | < EQL | | | | | ug/L |
| | 1,2-Dibromo-3-chloropropane | <10 | U | | < EQL | <10 | U | | < EQL | ug/L |
| | 1,2-Dibromoethane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,2-Dichlorobenzene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,2-Dichloroethane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,2-Dichloropropane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,3,5-Trinitrobenzene | <10 | JU | Q | < EQL | | | | | ug/L |
| | 1,3-Dichlorobenzene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,3-Dichloropropane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,3-Dinitrobenzene | <50 | JU | Q | < EQL | | | | | ug/L |
| | 1,4-Dichlorobenzene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,4-Naphthoquinone | <11 | JU | Q | < EQL | | | | | ug/L |
| | 1-Naphthylamine | <11 | JU | Q | < EQL | | | | | ug/L |
| | 2,2-Dichloropropane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 2,2-Oxybis(1-chloropropane) | <11 | JU | Q | < EQL | | | | | ug/L |
| | 2,3,4,6-Tetrachlorophenol | <11 | JU | Q | < EQL | | | | | ug/L |
| | 2,4,5-T | <2 | U | Q | < EQL | | | | | ug/L |
| | 2,4,5-TP (Silvex) | <2 | U | Q | < EQL | | | | | ug/L |
| | 2,4,5-Trichlorophenol | <11 | JU | Q | < EQL | | | | | ug/L |
| | 2,4,6-Trichlorophenol | <26 | JU | Q | < EQL | | | | | ug/L |
| | 2,4-Dichlorophenol | <11 | JU | Q | < EQL | | | | | ug/L |
| | 2,4-Dichlorophenoxyacetic acid | <2 | U | Q | < EQL | | | | | ug/L |
| | 2,4-Dimethyl phenol | <11 | JU | Q | < EQL | | | | | ug/L |
| | 2,4-Dinitrophenol | <26 | JU | Q | < EQL | | | | | ug/L |
| | 2,4-Dinitrotoluene | <11 | JU | Q | < EQL | | | | | ug/L |
| | 2,6-Dichlorophenol | <25 | JU | Q | < EQL | | | | | ug/L |
| | 2,6-Dinitrotoluene | <11 | JU | Q | < EQL | | | | | ug/L |
| | 2-Acetylaminofluorene | <11 | JU | Q | < EQL | | | | | ug/L |
| | 2-Chloronaphthalene | <11 | JU | Q | < EQL | | | | | ug/L |
| | 2-Chlorophenol | <11 | JU | Q | < EQL | | | | | ug/L |
| | 2-Hexanone | <20 | U | | < EQL | <20 | U | | < EQL | ug/L |
| | 2-Methyl-4,6-dinitrophenol | <26 | JU | Q | < EQL | | | | | ug/L |
| | 2-Methylnaphthalene | <11 | JU | Q | < EQL | | | | | ug/L |
| | 2-Naphthylamine | <11 | JU | Q | < EQL | | | | | ug/L |
| | 2-Nitrophenol | <11 | JU | Q | < EQL | | | | | ug/L |
| | 2-sec-Butyl-4,6-dinitrophenol | <2 | U | Q | < EQL | | | | | ug/L |
| | 3,3"-Dichlorobenzidine | <11 | JU | Q | < EQL | | | | | ug/L |
| | 3,3"-Dimethylbenzidine | <21 | JU | Q | < EQL | | | | | ug/L |
| | 3-Methylcholanthrene | <11 | JU | Q | < EQL | | | | | ug/L |
| | 4-Aminobiphenyl | <11 | JU | Q | < EQL | | | | | ug/L |
| | 4-Bromophenyl phenyl ether | <11 | JU | Q | < EQL | | | | | ug/L |
| | 4-Chloro-m-cresol | <11 | JU | Q | < EQL | | | | | ug/L |

Notes: Concentrations in bold italics exceed the groundwater protection standards listed in Appendix A. Bold italics were not assigned if the result was qualified with a "J", "R", "L", or "U" qualifier.

+ = exceeded the Groundwater Protection Standards listed in Appendix A.

Table B. Assessment Monitoring Results for Groundwater Wells (Cont.)

WELL LFW 75C

ANALYTICAL DATA

Groundwater Protection Standard

| ST | Parameter | 1st Half | CLP | EPA | Filt. | 2nd Half | CLP | EPA | Filt. | Unit |
|----|--------------------------------------|----------|-----|-----|----------|----------|-----|-----|-------|------|
| | 4-Chloroaniline | <11 | JU | Q | < EQL | | | | | ug/L |
| | 4-Chlorophenyl phenyl ether | <11 | JU | Q | < EQL | | | | | ug/L |
| | 4-Nitrophenol | <26 | JU | Q | < EQL | | | | | ug/L |
| | 5-Nitro-o-toluidine | <11 | JU | Q | < EQL | | | | | ug/L |
| | 7,12-Dimethylbenz(a)anthracene | <11 | JU | Q | < EQL | | | | | ug/L |
| | Acenaphthene | <11 | JU | Q | < EQL | | | | | ug/L |
| | Acenaphthylene | <11 | JU | Q | < EQL | | | | | ug/L |
| | Acetone | <20 | U | | < EQL | <20 | U | | < EQL | ug/L |
| | Acetonitrile (Methyl cyanide) | <200 | U | | < EQL | <200 | U | | < EQL | ug/L |
| | Acetophenone | <11 | JU | Q | < EQL | | | | | ug/L |
| | Acrolein | <50 | U | | < EQL | <50 | U | | < EQL | ug/L |
| | Acrylonitrile | <10 | U | | < EQL | <10 | U | | < EQL | ug/L |
| | Aldrin | <.1 | U | | < EQL | | | | | ug/L |
| | Allyl chloride | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Anthracene | <11 | JU | Q | < EQL | | | | | ug/L |
| | Antimony, total recoverable | <100 | U | | < EQL | <100 | U | | < EQL | ug/L |
| | Arsenic, total recoverable | <10 | U | | < EQL | <3.41 | JU | | < EQL | ug/L |
| | Barium, total recoverable | 6.74 | J | I | NDD | 6.94 | J | I | NDD | ug/L |
| | Benzene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Benzo(a)anthracene | <11 | JU | Q | < EQL | | | | | ug/L |
| | Benzo(a)pyrene | <11 | JU | Q | < EQL | | | | | ug/L |
| | Benzo(b)fluoranthene | <11 | JU | Q | < EQL | | | | | ug/L |
| | Benzo(g,h,i)perylene | <11 | JU | Q | < EQL | | | | | ug/L |
| | Benzo(k)fluoranthene | <11 | JU | Q | < EQL | | | | | ug/L |
| | Benzyl alcohol | | | | | | | | | |
| | Beryllium, total recoverable | <.16 | JU | | < EQL | <.32 | JU | | < EQL | ug/L |
| | Bis(2-chloroethoxy) methane | <11 | JU | Q | < EQL | | | | | ug/L |
| | Bis(2-chloroethyl) ether | <11 | JU | Q | < EQL | | | | | ug/L |
| | Bis(2-ethylhexyl) phthalate | <11 | JU | Q | < EQL | | | | | ug/L |
| | Bromochloromethane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Bromodichloromethane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Bromoform | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Bromomethane (Methyl bromide) | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Butylbenzyl phthalate | <11 | JU | Q | < EQL | | | | | ug/L |
| | Cadmium, total recoverable | <10 | U | | < EQL | <10 | U | | < EQL | ug/L |
| | Carbon disulfide | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Carbon tetrachloride | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Chlorobenzene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Chlorobenzilate | <11 | JU | Q | < EQL | | | | | ug/L |
| | Chloroethane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Chloroethene (Vinyl chloride) | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Chloroform | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Chloromethane (Methyl chloride) | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Chloroprene | <20 | U | | < EQL | <20 | U | | < EQL | ug/L |
| | Chromium, total recoverable | <10 | U | | < EQL | <10 | U | | < EQL | ug/L |
| | Chrysene | <11 | JU | Q | < EQL | | | | | ug/L |
| | Cobalt, total recoverable | <20 | U | | < EQL | 8.34 | J | I | NDD | ug/L |
| | Copper, total recoverable | <2.17 | JU | | < EQL | <1.9 | JU | | < EQL | ug/L |
| | Cyanide | 10 | R | Q | Rejected | | | | | ug/L |
| | Di-n-butyl phthalate | <11 | JU | Q | < EQL | <9.7 | U | | < EQL | ug/L |
| | Di-n-octyl phthalate | <11 | JU | Q | < EQL | | | | | ug/L |
| | Diallate | <11 | JU | Q | < EQL | | | | | ug/L |
| | Dibenz(a,h)anthracene | <11 | JU | Q | < EQL | | | | | ug/L |
| | Dibenzofuran | <11 | JU | Q | < EQL | | | | | ug/L |
| | Dibromochloromethane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Dibromomethane (Methylene bromide) | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Dichlorodifluoromethane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Dichloromethane (Methylene chloride) | 1.9 | J | I | NDD | <10 | U | | < EQL | ug/L |
| | Dieldrin | <.21 | U | | < EQL | | | | | ug/L |
| | Diethyl phthalate | <11 | JU | Q | < EQL | | | | | ug/L |
| | Dimethoate | <11 | JU | Q | < EQL | | | | | ug/L |
| | Dimethyl phthalate | <11 | JU | Q | < EQL | | | | | ug/L |
| | Diphenylamine | <11 | JU | Q | < EQL | | | | | ug/L |
| | Disulfoton | <11 | JU | Q | < EQL | | | | | ug/L |
| | Endosulfan I | <.1 | U | | < EQL | | | | | ug/L |
| | Endosulfan II | <.21 | U | | < EQL | | | | | ug/L |
| | Endosulfan sulfate | <.21 | U | | < EQL | | | | | ug/L |
| | Endrin | <.21 | U | | < EQL | | | | | ug/L |

Notes: Concentrations in bold italics exceed the groundwater protection standards listed in Appendix A. Bold italics were not assigned if the result was qualified with a "J", "R", "L", or "U" qualifier.

+ = exceeded the Groundwater Protection Standards listed in Appendix A.

Table B. Assessment Monitoring Results for Groundwater Wells (Cont.)

WELL LFW 75C

ANALYTICAL DATA

Groundwater Protection Standard

| ST | Parameter | 1st Half | CLP | EPA | Filt. | 2nd Half | CLP | EPA | Filt. | Unit |
|----|---------------------------------|----------|-----|-----|----------|----------|-----|-----|-------|-------|
| | Endrin aldehyde | <.21 | U | | < EQL | | | | | ug/L |
| | Ethyl methacrylate | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Ethyl methanesulfonate | <11 | JU | Q | < EQL | | | | | ug/L |
| | Ethylbenzene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Famphur | <200 | JU | Q | < EQL | | | | | ug/L |
| | Fluoranthene | <11 | JU | Q | < EQL | | | | | ug/L |
| | Fluorene | <11 | JU | Q | < EQL | | | | | ug/L |
| | Heptachlor | <.1 | U | | < EQL | | | | | ug/L |
| | Heptachlor epoxide | <.1 | U | | < EQL | | | | | ug/L |
| | Hexachlorobenzene | <11 | JU | Q | < EQL | | | | | ug/L |
| | Hexachlorobutadiene | <11 | JU | Q | < EQL | | | | | ug/L |
| | Hexachlorocyclopentadiene | <11 | JU | Q | < EQL | | | | | ug/L |
| | Hexachloroethane | <11 | JU | Q | < EQL | | | | | ug/L |
| | Hexachloropropene | <53 | JU | Q | < EQL | | | | | ug/L |
| | Indeno(1,2,3-c,d)pyrene | <11 | JU | Q | < EQL | | | | | ug/L |
| | Iodomethane (Methyl iodide) | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Isobutyl alcohol | <500 | U | | < EQL | <500 | U | | < EQL | ug/L |
| | Isodrin | <11 | JU | Q | < EQL | | | | | ug/L |
| | Isophorone | <11 | JU | Q | < EQL | | | | | ug/L |
| | Isosafrole | <11 | JU | Q | < EQL | | | | | ug/L |
| | Kepone | <11 | JU | Q | < EQL | | | | | ug/L |
| | Lead, total recoverable | 3.44 | J | I | NDD | <10 | U | | < EQL | ug/L |
| | Lindane | <.1 | U | | < EQL | | | | | ug/L |
| | Mercury, total recoverable | <.5 | U | | < EQL | <.5 | U | | < EQL | ug/L |
| | Methacrylonitrile | <200 | U | | < EQL | <200 | U | | < EQL | ug/L |
| | Methapyrilene | <11 | JU | Q | < EQL | | | | | ug/L |
| | Methoxychlor | <1 | U | | < EQL | | | | | ug/L |
| | Methyl ethyl ketone | <20 | U | | < EQL | <20 | U | | < EQL | ug/L |
| | Methyl isobutyl ketone | <10 | U | | < EQL | <10 | U | | < EQL | ug/L |
| | Methyl methacrylate | <20 | U | | < EQL | <20 | U | | < EQL | ug/L |
| | Methyl methanesulfonate | <11 | JU | Q | < EQL | | | | | ug/L |
| | N-Nitrosodi-n-butylamine | <11 | JU | Q | < EQL | | | | | ug/L |
| | N-Nitrosodiethylamine | <11 | JU | Q | < EQL | | | | | ug/L |
| | N-Nitrosodimethylamine | <26 | JU | Q | < EQL | | | | | ug/L |
| | N-Nitrosodiphenylamine | <11 | JU | Q | < EQL | | | | | ug/L |
| | N-Nitrosodipropylamine | <11 | JU | Q | < EQL | | | | | ug/L |
| | N-Nitrosomethylethylamine | <11 | JU | Q | < EQL | | | | | ug/L |
| | N-Nitrosopiperidine | <11 | JU | Q | < EQL | | | | | ug/L |
| | N-Nitrosopyrrolidine | <11 | JU | Q | < EQL | | | | | ug/L |
| | Naphthalene | <11 | JU | Q | < EQL | | | | | ug/L |
| | Nickel, total recoverable | <50 | U | | < EQL | <50 | U | | < EQL | ug/L |
| | Nitrobenzene | <11 | JU | Q | < EQL | | | | | ug/L |
| | O,O,O-Triethyl phosphorothioate | <11 | JU | Q | < EQL | | | | | ug/L |
| | PCB | | | | | | | | | |
| | PCB 1016 | <1 | U | | < EQL | | | | | ug/L |
| | PCB 1232 | <1 | U | | < EQL | | | | | ug/L |
| | PCB 1242 | <2.1 | U | | < EQL | | | | | ug/L |
| | PCB 1248 | <1 | U | | < EQL | | | | | ug/L |
| | PCB 1254 | <1 | U | | < EQL | | | | | ug/L |
| | PCB 1260 | <1 | U | | < EQL | | | | | ug/L |
| | Parathion | <11 | JU | Q | < EQL | | | | | ug/L |
| | Parathion methyl | <11 | JU | Q | < EQL | | | | | ug/L |
| | Pentachlorobenzene | <53 | JU | Q | < EQL | | | | | ug/L |
| | Pentachloronitrobenzene | <11 | JU | Q | < EQL | | | | | ug/L |
| | Pentachlorophenol | <26 | JU | Q | < EQL | | | | | ug/L |
| | Phenacetin | <11 | JU | Q | < EQL | | | | | ug/L |
| | Phenanthrene | <11 | JU | Q | < EQL | | | | | ug/L |
| | Phenol | <11 | JU | Q | < EQL | | | | | ug/L |
| | Phorate | <11 | JU | Q | < EQL | | | | | ug/L |
| | Pronamid | <11 | JU | Q | < EQL | | | | | ug/L |
| | Propionitrile | <200 | U | | < EQL | <200 | U | | < EQL | ug/L |
| | Pyrene | <11 | JU | Q | < EQL | | | | | ug/L |
| | Safrole | <11 | JU | Q | < EQL | | | | | ug/L |
| | Selenium, total recoverable | <10 | U | | < EQL | <10 | U | | < EQL | ug/L |
| | Silver, total recoverable | <20 | U | | < EQL | <20 | U | | < EQL | ug/L |
| | Specific conductance | 29.6 | | | < 37 | 30.5 | | | < 37 | uS/cm |
| | Styrene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Sulfide | 1000 | R | Q | Rejected | | | | | ug/L |

Notes: Concentrations in bold italics exceed the groundwater protection standards listed in Appendix A. Bold italics were not assigned if the result was qualified with a "J", "R", "L", or "U" qualifier.

+ = exceeded the Groundwater Protection Standards listed in Appendix A.

Table B. Assessment Monitoring Results for Groundwater Wells (Cont.)

WELL LFW 75C

ANALYTICAL DATA

Groundwater Protection Standard

| ST | Parameter | 1st Half | CLP | EPA | Filt. | 2nd Half | CLP | EPA | Filt. | Unit |
|----|-----------------------------|----------|-----|-----|--------|----------|-----|-----|-------|------|
| | Tetrachloroethylene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Thallium, total recoverable | <5.04 | JU | | < EQL | <10 | U | | < EQL | ug/L |
| | Thionazin | <11 | JU | Q | < EQL | | | | | ug/L |
| | Tin, total recoverable | <200 | U | | < EQL | | | | | ug/L |
| | Toluene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Toxaphene | <2.1 | U | | < EQL | | | | | ug/L |
| | Trichloroethylene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Trichlorofluoromethane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Vanadium, total recoverable | <10 | U | | < EQL | <10 | U | | < EQL | ug/L |
| | Vinyl acetate | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Xylenes | <10 | U | | < EQL | <10 | U | | < EQL | ug/L |
| | Zinc, total recoverable | 20.2 | | | < 29.3 | <8.28 | JU | | < EQL | ug/L |
| | alpha-Benzene hexachloride | <.1 | U | | < EQL | | | | | ug/L |
| | alpha-Chlordane | <.1 | U | | < EQL | | | | | ug/L |
| | beta-Benzene hexachloride | <.1 | U | | < EQL | <.095 | U | | < EQL | ug/L |
| | cis-1,2-Dichloroethylene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | cis-1,3-Dichloropropene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | delta-Benzene hexachloride | <.1 | U | | < EQL | | | | | ug/L |
| | gamma-Chlordane | <.1 | U | | < EQL | | | | | ug/L |
| | m-Cresol (3-Methylphenol) | | | | | | | | | |
| | m-Nitroaniline | <26 | JU | Q | < EQL | | | | | ug/L |
| | o-Cresol (2-Methylphenol) | <11 | JU | Q | < EQL | | | | | ug/L |
| | o-Nitroaniline | <26 | JU | Q | < EQL | | | | | ug/L |
| | o-Toluidine | <11 | JU | Q | < EQL | | | | | ug/L |
| | p,p"-DDD | <.21 | U | | < EQL | | | | | ug/L |
| | p,p"-DDE | <.21 | U | | < EQL | | | | | ug/L |
| | p,p"-DDT | <.21 | U | | < EQL | | | | | ug/L |
| | p-Cresol (4-Methylphenol) | <10 | JU | Q | < EQL | | | | | ug/L |
| | p-Dimethylaminoazobenzene | <11 | JU | Q | < EQL | | | | | ug/L |
| | p-Nitroaniline | <11 | JU | Q | < EQL | | | | | ug/L |
| | p-Phenylenediamine | <11 | JU | Q | < EQL | | | | | ug/L |
| | pH | 5.5 | J | Q | NDD | 5.02 | J | Q | NDD | pH |
| | trans-1,2-Dichloroethylene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | trans-1,3-Dichloropropene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | trans-1,4-Dichloro-2-butene | <20 | U | | < EQL | <20 | U | | < EQL | ug/L |

Notes: Concentrations in bold italics exceed the groundwater protection standards listed in Appendix A. Bold italics were not assigned if the result was qualified with a "J", "R", "L", or "U" qualifier.

+ = exceeded the Groundwater Protection Standards listed in Appendix A.

Table B. Assessment Monitoring Results for Groundwater Wells (Cont.)

WELL LFW 75D

| SRS Coord. | Lat/Longitude | Screen Zone Elevation | Top of Casing | Casing | Pump | Screen Zone |
|----------------------|----------------------------------|-----------------------|---------------|---------|------|-------------|
| N 85868 E 45355.6 | 33.28968 Deg N -81.7142 Deg W | 166 - 151 ft msl | 198.2 ft msl | 4 " PVC | S | Upper |

SAMPLE DATE 03/08/00 10/02/00

FIELD DATA

| Parameter | 1st Half | 2nd Half | Unit |
|---------------------|----------|----------|-------------|
| Water Elevation | 161.01 | 160.5 | ft msl |
| pH | 5.1 | 5.1 | pH |
| Sp. Conductance | 25 | 26 | uS/cm |
| Water temperature | 20.7 | 21.3 | deg. C |
| Alkalinity as CaCO3 | 3 | 0 | mg/L |
| Turbidity | .5 | 1.4 | NTU |
| Volumes purged | 0 | 5.03 | well volume |
| Sampling code | | | |

ANALYTICAL DATA

Groundwater Protection Standard

| ST | Parameter | 1st Half | CLP | EPA | Filt. | 2nd Half | CLP | EPA | Filt. | Unit |
|----|--------------------------------|----------|-----|-----|-------|----------|-----|-----|-------|------|
| | 1,1,1,2-Tetrachloroethane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,1,1-Trichloroethane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,1,2,2-Tetrachloroethane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,1,2-Trichloroethane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,1-Dichloroethane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,1-Dichloroethylene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,1-Dichloropropene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,2,3-Trichloropropane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,2,4,5-Tetrachlorobenzene | <53 | JU | Q | < EQL | | | | | ug/L |
| | 1,2,4-Trichlorobenzene | <11 | JU | Q | < EQL | | | | | ug/L |
| | 1,2-Dibromo-3-chloropropane | <10 | U | | < EQL | <10 | U | | < EQL | ug/L |
| | 1,2-Dibromoethane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,2-Dichlorobenzene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,2-Dichloroethane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,2-Dichloropropane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,3,5-Trinitrobenzene | <10 | JU | Q | < EQL | | | | | ug/L |
| | 1,3-Dichlorobenzene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,3-Dichloropropane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,3-Dinitrobenzene | <50 | JU | Q | < EQL | | | | | ug/L |
| | 1,4-Dichlorobenzene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,4-Naphthoquinone | <11 | JU | Q | < EQL | | | | | ug/L |
| | 1-Naphthylamine | <11 | JU | Q | < EQL | | | | | ug/L |
| | 2,2-Dichloropropane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 2,2-Oxybis(1-chloropropane) | <11 | JU | Q | < EQL | | | | | ug/L |
| | 2,3,4,6-Tetrachlorophenol | <11 | JU | Q | < EQL | | | | | ug/L |
| | 2,4,5-T | <2 | U | Q | < EQL | | | | | ug/L |
| | 2,4,5-TP (Silvex) | <2 | U | Q | < EQL | | | | | ug/L |
| | 2,4,5-Trichlorophenol | <11 | JU | Q | < EQL | | | | | ug/L |
| | 2,4,6-Trichlorophenol | <26 | JU | Q | < EQL | | | | | ug/L |
| | 2,4-Dichlorophenol | <11 | JU | Q | < EQL | | | | | ug/L |
| | 2,4-Dichlorophenoxyacetic acid | <2 | U | Q | < EQL | | | | | ug/L |
| | 2,4-Dimethyl phenol | <11 | JU | Q | < EQL | | | | | ug/L |
| | 2,4-Dinitrophenol | <26 | JU | Q | < EQL | | | | | ug/L |
| | 2,4-Dinitrotoluene | <11 | JU | Q | < EQL | | | | | ug/L |
| | 2,6-Dichlorophenol | <25 | JU | Q | < EQL | | | | | ug/L |
| | 2,6-Dinitrotoluene | <11 | JU | Q | < EQL | | | | | ug/L |
| | 2-Acetylaminofluorene | <11 | JU | Q | < EQL | | | | | ug/L |
| | 2-Chloronaphthalene | <11 | JU | Q | < EQL | | | | | ug/L |
| | 2-Chlorophenol | <11 | JU | Q | < EQL | | | | | ug/L |
| | 2-Hexanone | <20 | U | | < EQL | <20 | U | | < EQL | ug/L |
| | 2-Methyl-4,6-dinitrophenol | <26 | JU | Q | < EQL | | | | | ug/L |
| | 2-Methylnaphthalene | <11 | JU | Q | < EQL | | | | | ug/L |
| | 2-Naphthylamine | <11 | JU | Q | < EQL | | | | | ug/L |
| | 2-Nitrophenol | <11 | JU | Q | < EQL | | | | | ug/L |
| | 2-sec-Butyl-4,6-dinitrophenol | <2 | U | Q | < EQL | | | | | ug/L |
| | 3,3'-Dichlorobenzidine | <11 | JU | Q | < EQL | | | | | ug/L |
| | 3,3'-Dimethylbenzidine | <21 | JU | Q | < EQL | | | | | ug/L |
| | 3-Methylcholanthrene | <11 | JU | Q | < EQL | | | | | ug/L |
| | 4-Aminobiphenyl | <11 | JU | Q | < EQL | | | | | ug/L |
| | 4-Bromophenyl phenyl ether | <11 | JU | Q | < EQL | | | | | ug/L |
| | 4-Chloro-m-cresol | <11 | JU | Q | < EQL | | | | | ug/L |

Notes: Concentrations in bold italics exceed the groundwater protection standards listed in Appendix A. Bold italics were not assigned if the result was qualified with a "J", "R", "L", or "U" qualifier.

+ = exceeded the Groundwater Protection Standards listed in Appendix A.

Table B. Assessment Monitoring Results for Groundwater Wells (Cont.)

WELL LFW 75D

ANALYTICAL DATA

Groundwater Protection Standard

| ST | Parameter | 1st Half | CLP | EPA | Filt. | 2nd Half | CLP | EPA | Filt. | Unit |
|----|--------------------------------------|----------|-----|-----|----------|----------|-----|-----|--------|------|
| | 4-Chloroaniline | <11 | JU | Q | < EQL | | | | | ug/L |
| | 4-Chlorophenyl phenyl ether | <11 | JU | Q | < EQL | | | | | ug/L |
| | 4-Nitrophenol | <26 | JU | Q | < EQL | | | | | ug/L |
| | 5-Nitro-o-toluidine | <11 | JU | Q | < EQL | | | | | ug/L |
| | 7,12-Dimethylbenz(a)anthracene | <11 | JU | Q | < EQL | | | | | ug/L |
| | Acenaphthene | <11 | JU | Q | < EQL | | | | | ug/L |
| | Acenaphthylene | <11 | JU | Q | < EQL | | | | | ug/L |
| | Acetone | <20 | U | | < EQL | <20 | U | | < EQL | ug/L |
| | Acetonitrile (Methyl cyanide) | <200 | U | | < EQL | <200 | U | | < EQL | ug/L |
| | Acetophenone | <11 | JU | Q | < EQL | | | | | ug/L |
| | Acrolein | <50 | U | | < EQL | <50 | U | | < EQL | ug/L |
| | Acrylonitrile | <10 | U | | < EQL | <10 | U | | < EQL | ug/L |
| | Aldrin | <.1 | U | | < EQL | | | | | ug/L |
| | Allyl chloride | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Anthracene | <11 | JU | Q | < EQL | | | | | ug/L |
| | Antimony, total recoverable | <100 | U | | < EQL | <100 | U | | < EQL | ug/L |
| | Arsenic, total recoverable | <10 | U | | < EQL | <10 | U | | < EQL | ug/L |
| | Barium, total recoverable | 11.6 | | | < 2000 | 11.9 | | | < 2000 | ug/L |
| | Benzene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Benzo(a)anthracene | <11 | JU | Q | < EQL | | | | | ug/L |
| | Benzo(a)pyrene | <11 | JU | Q | < EQL | | | | | ug/L |
| | Benzo(b)fluoranthene | <11 | JU | Q | < EQL | | | | | ug/L |
| | Benzo(g,h,i)perylene | <11 | JU | Q | < EQL | | | | | ug/L |
| | Benzo(k)fluoranthene | <11 | JU | Q | < EQL | | | | | ug/L |
| | Benzyl alcohol | | | | | | | | | |
| | Beryllium, total recoverable | <1 | U | | < EQL | <.39 | JU | | < EQL | ug/L |
| | Bis(2-chloroethoxy) methane | <11 | JU | Q | < EQL | | | | | ug/L |
| | Bis(2-chloroethyl) ether | <11 | JU | Q | < EQL | | | | | ug/L |
| | Bis(2-ethylhexyl) phthalate | <11 | JU | Q | < EQL | | | | | ug/L |
| | Bromochloromethane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Bromodichloromethane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Bromoform | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Bromomethane (Methyl bromide) | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Butylbenzyl phthalate | <11 | JU | Q | < EQL | | | | | ug/L |
| | Cadmium, total recoverable | <10 | U | | < EQL | <10 | U | | < EQL | ug/L |
| | Carbon disulfide | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Carbon tetrachloride | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Chlorobenzene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Chlorobenzilate | <11 | JU | Q | < EQL | | | | | ug/L |
| | Chloroethane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Chloroethene (Vinyl chloride) | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Chloroform | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Chloromethane (Methyl chloride) | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Chloroprene | <20 | U | | < EQL | <20 | U | | < EQL | ug/L |
| | Chromium, total recoverable | <10 | U | | < EQL | <10 | U | | < EQL | ug/L |
| | Chrysene | <11 | JU | Q | < EQL | | | | | ug/L |
| | Cobalt, total recoverable | <20 | U | | < EQL | <20 | U | | < EQL | ug/L |
| | Copper, total recoverable | <20 | U | | < EQL | <1.9 | JU | | < EQL | ug/L |
| | Cyanide | 10 | R | Q | Rejected | | | | | ug/L |
| | Di-n-butyl phthalate | <11 | JU | Q | < EQL | <9.8 | U | | < EQL | ug/L |
| | Di-n-octyl phthalate | <11 | JU | Q | < EQL | | | | | ug/L |
| | Diallate | <11 | JU | Q | < EQL | | | | | ug/L |
| | Dibenz(a,h)anthracene | <11 | JU | Q | < EQL | | | | | ug/L |
| | Dibenzofuran | <11 | JU | Q | < EQL | | | | | ug/L |
| | Dibromochloromethane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Dibromomethane (Methylene bromide) | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Dichlorodifluoromethane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Dichloromethane (Methylene chloride) | 1.8 | J | I | NDD | <10 | U | | < EQL | ug/L |
| | Dieldrin | <.21 | U | | < EQL | | | | | ug/L |
| | Diethyl phthalate | <11 | JU | Q | < EQL | | | | | ug/L |
| | Dimethoate | <11 | JU | Q | < EQL | | | | | ug/L |
| | Dimethyl phthalate | <11 | JU | Q | < EQL | | | | | ug/L |
| | Diphenylamine | <11 | JU | Q | < EQL | | | | | ug/L |
| | Disulfoton | <11 | JU | Q | < EQL | | | | | ug/L |
| | Endosulfan I | <.1 | U | | < EQL | | | | | ug/L |
| | Endosulfan II | <.21 | U | | < EQL | | | | | ug/L |
| | Endosulfan sulfate | <.21 | U | | < EQL | | | | | ug/L |
| | Endrin | <.21 | U | | < EQL | | | | | ug/L |

Notes: Concentrations in bold italics exceed the groundwater protection standards listed in Appendix A. Bold italics were not assigned if the result was qualified with a "J", "R", "L", or "U" qualifier.

+ = exceeded the Groundwater Protection Standards listed in Appendix A.

Table B. Assessment Monitoring Results for Groundwater Wells (Cont.)

WELL LFW 75D

ANALYTICAL DATA

Groundwater Protection Standard

| ST | Parameter | 1st Half | CLP | EPA | Filt. | 2nd Half | CLP | EPA | Filt. | Unit |
|----|---------------------------------|----------|-----|-----|----------|----------|-----|-----|-------|-------|
| | Endrin aldehyde | <.21 | U | | < EQL | | | | | ug/L |
| | Ethyl methacrylate | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Ethyl methanesulfonate | <11 | JU | Q | < EQL | | | | | ug/L |
| | Ethylbenzene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Famphur | <200 | JU | Q | < EQL | | | | | ug/L |
| | Fluoranthene | <11 | JU | Q | < EQL | | | | | ug/L |
| | Fluorene | <11 | JU | Q | < EQL | | | | | ug/L |
| | Heptachlor | <.1 | U | | < EQL | | | | | ug/L |
| | Heptachlor epoxide | <.1 | U | | < EQL | | | | | ug/L |
| | Hexachlorobenzene | <11 | JU | Q | < EQL | | | | | ug/L |
| | Hexachlorobutadiene | <11 | JU | Q | < EQL | | | | | ug/L |
| | Hexachlorocyclopentadiene | <11 | JU | Q | < EQL | | | | | ug/L |
| | Hexachloroethane | <11 | JU | Q | < EQL | | | | | ug/L |
| | Hexachloropropene | <53 | JU | Q | < EQL | | | | | ug/L |
| | Indeno(1,2,3-c,d)pyrene | <11 | JU | Q | < EQL | | | | | ug/L |
| | Iodomethane (Methyl iodide) | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Isobutyl alcohol | <500 | U | | < EQL | <500 | U | | < EQL | ug/L |
| | Isodrin | <11 | JU | Q | < EQL | | | | | ug/L |
| | Isophorone | <11 | JU | Q | < EQL | | | | | ug/L |
| | Isosafrole | <11 | JU | Q | < EQL | | | | | ug/L |
| | Kepone | <11 | JU | Q | < EQL | | | | | ug/L |
| | Lead, total recoverable | 3.74 | J | I | NDD | <1.3 | JU | | < EQL | ug/L |
| | Lindane | <.1 | U | | < EQL | | | | | ug/L |
| | Mercury, total recoverable | <.5 | U | | < EQL | <.5 | U | | < EQL | ug/L |
| | Methacrylonitrile | <200 | U | | < EQL | <200 | U | | < EQL | ug/L |
| | Methapyrilene | <11 | JU | Q | < EQL | | | | | ug/L |
| | Methoxychlor | <.1 | U | | < EQL | | | | | ug/L |
| | Methyl ethyl ketone | <20 | U | | < EQL | <20 | U | | < EQL | ug/L |
| | Methyl isobutyl ketone | <10 | U | | < EQL | <10 | U | | < EQL | ug/L |
| | Methyl methacrylate | <20 | U | | < EQL | <20 | U | | < EQL | ug/L |
| | Methyl methanesulfonate | <11 | JU | Q | < EQL | | | | | ug/L |
| | N-Nitrosodi-n-butylamine | <11 | JU | Q | < EQL | | | | | ug/L |
| | N-Nitrosodiethylamine | <11 | JU | Q | < EQL | | | | | ug/L |
| | N-Nitrosodimethylamine | <26 | JU | Q | < EQL | | | | | ug/L |
| | N-Nitrosodiphenylamine | <11 | JU | Q | < EQL | | | | | ug/L |
| | N-Nitrosodipropylamine | <11 | JU | Q | < EQL | | | | | ug/L |
| | N-Nitrosomethylethylamine | <11 | JU | Q | < EQL | | | | | ug/L |
| | N-Nitrosopiperidine | <11 | JU | Q | < EQL | | | | | ug/L |
| | N-Nitrosopyrrolidine | <11 | JU | Q | < EQL | | | | | ug/L |
| | Naphthalene | <11 | JU | Q | < EQL | | | | | ug/L |
| | Nickel, total recoverable | <50 | U | | < EQL | <50 | U | | < EQL | ug/L |
| | Nitrobenzene | <11 | JU | Q | < EQL | | | | | ug/L |
| | O,O,O-Triethyl phosphorothioate | <11 | JU | Q | < EQL | | | | | ug/L |
| | PCB | | | | | | | | | |
| | PCB 1016 | <.1 | U | | < EQL | | | | | ug/L |
| | PCB 1232 | <.1 | U | | < EQL | | | | | ug/L |
| | PCB 1242 | <2.1 | U | | < EQL | | | | | ug/L |
| | PCB 1248 | <.1 | U | | < EQL | | | | | ug/L |
| | PCB 1254 | <.1 | U | | < EQL | | | | | ug/L |
| | PCB 1260 | <.1 | U | | < EQL | | | | | ug/L |
| | Parathion | <11 | JU | Q | < EQL | | | | | ug/L |
| | Parathion methyl | <11 | JU | Q | < EQL | | | | | ug/L |
| | Pentachlorobenzene | <53 | JU | Q | < EQL | | | | | ug/L |
| | Pentachloronitrobenzene | <11 | JU | Q | < EQL | | | | | ug/L |
| | Pentachlorophenol | <26 | JU | Q | < EQL | | | | | ug/L |
| | Phenacetin | <11 | JU | Q | < EQL | | | | | ug/L |
| | Phenanthrene | <11 | JU | Q | < EQL | | | | | ug/L |
| | Phenol | <11 | JU | Q | < EQL | | | | | ug/L |
| | Phorate | <11 | JU | Q | < EQL | | | | | ug/L |
| | Pronamid | <11 | JU | Q | < EQL | | | | | ug/L |
| | Propionitrile | <200 | U | | < EQL | <200 | U | | < EQL | ug/L |
| | Pyrene | <11 | JU | Q | < EQL | | | | | ug/L |
| | Safrole | <11 | JU | Q | < EQL | | | | | ug/L |
| | Selenium, total recoverable | <10 | U | | < EQL | <10 | U | | < EQL | ug/L |
| | Silver, total recoverable | <20 | U | | < EQL | <3.33 | JU | | < EQL | ug/L |
| | Specific conductance | 24 | | | < 37 | 30.5 | | | < 37 | uS/cm |
| | Styrene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Sulfide | 1000 | R | Q | Rejected | | | | | ug/L |

Notes: Concentrations in bold italics exceed the groundwater protection standards listed in Appendix A. Bold italics were not assigned if the result was qualified with a "J", "R", "L", or "U" qualifier.

+ = exceeded the Groundwater Protection Standards listed in Appendix A.

Table B. Assessment Monitoring Results for Groundwater Wells (Cont.)

WELL LFW 75D

ANALYTICAL DATA

Groundwater Protection Standard

| ST | Parameter | 1st Half | CLP | EPA | Filt. | 2nd Half | CLP | EPA | Filt. | Unit |
|----|-----------------------------|----------|-----|-----|--------|----------|-----|-----|--------|------|
| | Tetrachloroethylene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Thallium, total recoverable | <10 | U | | < EQL | <10 | U | | < EQL | ug/L |
| | Thionazin | <11 | JU | Q | < EQL | | | | | ug/L |
| | Tin, total recoverable | <200 | U | | < EQL | | | | | ug/L |
| | Toluene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Toxaphene | <2.1 | U | | < EQL | | | | | ug/L |
| | Trichloroethylene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Trichlorofluoromethane | 13 | | | < 20 | 21 | J | K | NDD | ug/L |
| | Vanadium, total recoverable | <10 | U | | < EQL | <10 | U | | < EQL | ug/L |
| | Vinyl acetate | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Xylenes | <10 | U | | < EQL | <10 | U | | < EQL | ug/L |
| | Zinc, total recoverable | 28 | | | < 29.3 | 21.5 | | | < 29.3 | ug/L |
| | alpha-Benzene hexachloride | <.1 | U | | < EQL | | | | | ug/L |
| | alpha-Chlordane | <.1 | U | | < EQL | | | | | ug/L |
| | beta-Benzene hexachloride | <.1 | U | | < EQL | <.095 | U | | < EQL | ug/L |
| | cis-1,2-Dichloroethylene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | cis-1,3-Dichloropropene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | delta-Benzene hexachloride | <.1 | U | | < EQL | | | | | ug/L |
| | gamma-Chlordane | <.1 | U | | < EQL | | | | | ug/L |
| | m-Cresol (3-Methylphenol) | | | | | | | | | |
| | m-Nitroaniline | <26 | JU | Q | < EQL | | | | | ug/L |
| | o-Cresol (2-Methylphenol) | <11 | JU | Q | < EQL | | | | | ug/L |
| | o-Nitroaniline | <26 | JU | Q | < EQL | | | | | ug/L |
| | o-Toluidine | <11 | JU | Q | < EQL | | | | | ug/L |
| | p,p"-DDD | <.21 | U | | < EQL | | | | | ug/L |
| | p,p"-DDE | <.21 | U | | < EQL | | | | | ug/L |
| | p,p"-DDT | <.21 | U | | < EQL | | | | | ug/L |
| | p-Cresol (4-Methylphenol) | <10 | JU | Q | < EQL | | | | | ug/L |
| | p-Dimethylaminoazobenzene | <11 | JU | Q | < EQL | | | | | ug/L |
| | p-Nitroaniline | <11 | JU | Q | < EQL | | | | | ug/L |
| | p-Phenylenediamine | <11 | JU | Q | < EQL | | | | | ug/L |
| | pH | 5.81 | J | Q | NDD | 5.36 | J | Q | NDD | pH |
| | trans-1,2-Dichloroethylene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | trans-1,3-Dichloropropene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | trans-1,4-Dichloro-2-butene | <20 | U | | < EQL | <20 | U | | < EQL | ug/L |

Notes: Concentrations in bold italics exceed the groundwater protection standards listed in Appendix A. Bold italics were not assigned if the result was qualified with a "J", "R", "L", or "U" qualifier.

+ = exceeded the Groundwater Protection Standards listed in Appendix A.

Table B. Assessment Monitoring Results for Groundwater Wells (Cont.)

WELL LFW 76

| <u>SRS Coord.</u> | <u>Lat/Longitude</u> | <u>Screen Zone Elevation</u> | <u>Top of Casing</u> | <u>Casing</u> | <u>Pump</u> | <u>Screen Zone</u> |
|------------------------|----------------------------------|------------------------------|----------------------|---------------|-------------|--------------------|
| N 85682.1 E 44758.6 | 33.28830 Deg N -81.7154 Deg W | 157.9 - 142.9 ft msl | 221.9 ft msl | 2 " PVC | V | Upper |

SAMPLE DATE 02/28/00 10/02/00

FIELD DATA

| <u>Parameter</u> | <u>1st Half</u> | <u>2nd Half</u> | <u>Unit</u> |
|---------------------|-----------------|-----------------|-------------|
| Water Elevation | 157.69 | 157.13 | ft msl |
| pH | 5 | 5.4 | pH |
| Sp. Conductance | 25 | 23 | uS/cm |
| Water temperature | 22.7 | 35 | deg. C |
| Alkalinity as CaCO3 | 2 | 0 | mg/L |
| Turbidity | 14 | 25.2 | NTU |
| Volumes purged | 0 | 7.34 | well volume |
| Sampling code | NX | NX | |

ANALYTICAL DATA

Groundwater Protection Standard

| <u>ST</u> | <u>Parameter</u> | <u>1st Half</u> | <u>CLP</u> | <u>EPA</u> | <u>Filt.</u> | <u>2nd Half</u> | <u>CLP</u> | <u>EPA</u> | <u>Filt.</u> | <u>Unit</u> |
|-----------|--------------------------------|-----------------|------------|------------|--------------|-----------------|------------|------------|--------------|-------------|
| | 1,1,1,2-Tetrachloroethane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,1,1-Trichloroethane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,1,2,2-Tetrachloroethane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,1,2-Trichloroethane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,1-Dichloroethane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,1-Dichloroethylene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,1-Dichloropropene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,2,3-Trichloropropane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,2,4,5-Tetrachlorobenzene | <52 | U | | < EQL | | | | | ug/L |
| | 1,2,4-Trichlorobenzene | <11 | U | | < EQL | | | | | ug/L |
| | 1,2-Dibromo-3-chloropropane | <10 | U | | < EQL | <10 | U | | < EQL | ug/L |
| | 1,2-Dibromoethane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,2-Dichlorobenzene | <11 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,2-Dichloroethane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,2-Dichloropropane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,3,5-Trinitrobenzene | <10 | U | | < EQL | | | | | ug/L |
| | 1,3-Dichlorobenzene | <11 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,3-Dichloropropane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,3-Dinitrobenzene | <50 | U | | < EQL | | | | | ug/L |
| | 1,4-Dichlorobenzene | <11 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 1,4-Naphthoquinone | <11 | U | | < EQL | | | | | ug/L |
| | 1-Naphthylamine | <11 | U | | < EQL | | | | | ug/L |
| | 2,2-Dichloropropane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | 2,2-Oxybis(1-chloropropane) | <11 | U | | < EQL | | | | | ug/L |
| | 2,3,4,6-Tetrachlorophenol | <11 | U | | < EQL | | | | | ug/L |
| | 2,4,5-T | <2 | U | | < EQL | | | | | ug/L |
| | 2,4,5-TP (Silvex) | <2 | U | | < EQL | | | | | ug/L |
| | 2,4,5-Trichlorophenol | <11 | U | | < EQL | | | | | ug/L |
| | 2,4,6-Trichlorophenol | <26 | U | | < EQL | | | | | ug/L |
| | 2,4-Dichlorophenol | <11 | U | | < EQL | | | | | ug/L |
| | 2,4-Dichlorophenoxyacetic acid | <2 | U | | < EQL | | | | | ug/L |
| | 2,4-Dimethyl phenol | <11 | U | | < EQL | | | | | ug/L |
| | 2,4-Dinitrophenol | <26 | U | | < EQL | | | | | ug/L |
| | 2,4-Dinitrotoluene | <11 | U | | < EQL | | | | | ug/L |
| | 2,6-Dichlorophenol | <25 | U | | < EQL | | | | | ug/L |
| | 2,6-Dinitrotoluene | <11 | U | | < EQL | | | | | ug/L |
| | 2-Acetylaminofluorene | <11 | U | | < EQL | | | | | ug/L |
| | 2-Chloronaphthalene | <11 | U | | < EQL | | | | | ug/L |
| | 2-Chlorophenol | <11 | U | | < EQL | | | | | ug/L |
| | 2-Hexanone | <20 | U | | < EQL | <20 | U | | < EQL | ug/L |
| | 2-Methyl-4,6-dinitrophenol | <26 | U | | < EQL | | | | | ug/L |
| | 2-Methylnaphthalene | <11 | U | | < EQL | | | | | ug/L |
| | 2-Naphthylamine | <11 | U | | < EQL | | | | | ug/L |
| | 2-Nitrophenol | <11 | U | | < EQL | | | | | ug/L |
| | 2-sec-Butyl-4,6-dinitrophenol | <2 | JU | L | < EQL | | | | | ug/L |
| | 3,3"-Dichlorobenzidine | <11 | U | | < EQL | | | | | ug/L |
| | 3,3"-Dimethylbenzidine | <21 | U | | < EQL | | | | | ug/L |
| | 3-Methylcholanthrene | <11 | U | | < EQL | | | | | ug/L |
| | 4-Aminobiphenyl | <11 | U | | < EQL | | | | | ug/L |
| | 4-Bromophenyl phenyl ether | <11 | U | | < EQL | | | | | ug/L |
| | 4-Chloro-m-cresol | <11 | U | | < EQL | | | | | ug/L |

Notes: Concentrations in bold italics exceed the groundwater protection standards listed in Appendix A. Bold italics were not assigned if the result was qualified with a "J", "R", "L", or "U" qualifier.

+ = exceeded the Groundwater Protection Standards listed in Appendix A.

Table B. Assessment Monitoring Results for Groundwater Wells (Cont.)

WELL LFW 76

ANALYTICAL DATA

Groundwater Protection Standard

| ST | Parameter | 1st Half | CLP | EPA | Filt. | 2nd Half | CLP | EPA | Filt. | Unit |
|----|--------------------------------------|----------|-----|-----|-------|----------|-----|-----|-------|------|
| | 4-Chloroaniline | <11 | U | | < EQL | | | | | ug/L |
| | 4-Chlorophenyl phenyl ether | <11 | U | | < EQL | | | | | ug/L |
| | 4-Nitrophenol | <26 | U | | < EQL | | | | | ug/L |
| | 5-Nitro-o-toluidine | <11 | U | | < EQL | | | | | ug/L |
| | 7,12-Dimethylbenz(a)anthracene | <11 | U | | < EQL | | | | | ug/L |
| | Acenaphthene | <11 | U | | < EQL | | | | | ug/L |
| | Acenaphthylene | <11 | U | | < EQL | | | | | ug/L |
| | Acetone | <20 | U | | < EQL | <20 | U | | < EQL | ug/L |
| | Acetonitrile (Methyl cyanide) | <200 | U | | < EQL | <200 | U | | < EQL | ug/L |
| | Acetophenone | <11 | U | | < EQL | | | | | ug/L |
| | Acrolein | <50 | U | | < EQL | <50 | U | | < EQL | ug/L |
| | Acrylonitrile | <10 | U | | < EQL | <10 | U | | < EQL | ug/L |
| | Aldrin | <.1 | U | | < EQL | | | | | ug/L |
| | Allyl chloride | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Anthracene | <11 | U | | < EQL | | | | | ug/L |
| | Antimony, total recoverable | <100 | U | | < EQL | <100 | U | | < EQL | ug/L |
| | Arsenic, total recoverable | <10 | U | | < EQL | <10 | U | | < EQL | ug/L |
| | Barium, total recoverable | 12.7 | | | <2000 | 10 | J | I | NDD | ug/L |
| | Benzene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Benzo(a)anthracene | <11 | U | | < EQL | | | | | ug/L |
| | Benzo(a)pyrene | <11 | U | | < EQL | | | | | ug/L |
| | Benzo(b)fluoranthene | <11 | U | | < EQL | | | | | ug/L |
| | Benzo(g,h,i)perylene | <11 | U | | < EQL | | | | | ug/L |
| | Benzo(k)fluoranthene | <11 | U | | < EQL | | | | | ug/L |
| | Benzyl alcohol | | | | | | | | | |
| | Beryllium, total recoverable | <1 | U | | < EQL | <.11 | JU | | < EQL | ug/L |
| | Bis(2-chloroethoxy) methane | <11 | U | | < EQL | | | | | ug/L |
| | Bis(2-chloroethyl) ether | <11 | U | | < EQL | | | | | ug/L |
| | Bis(2-ethylhexyl) phthalate | <11 | U | | < EQL | | | | | ug/L |
| | Bromochloromethane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Bromodichloromethane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Bromoform | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Bromomethane (Methyl bromide) | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Butylbenzyl phthalate | <11 | U | | < EQL | | | | | ug/L |
| | Cadmium, total recoverable | <10 | U | | < EQL | <10 | U | | < EQL | ug/L |
| | Carbon disulfide | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Carbon tetrachloride | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Chlorobenzene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Chlorobenzilate | <11 | U | | < EQL | | | | | ug/L |
| | Chloroethane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Chloroethene (Vinyl chloride) | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Chloroform | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Chloromethane (Methyl chloride) | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Chloroprene | <20 | U | | < EQL | <20 | U | | < EQL | ug/L |
| | Chromium, total recoverable | <9.6 | U | V | < EQL | 25.8 | | | < 100 | ug/L |
| | Chrysene | <11 | U | | < EQL | | | | | ug/L |
| | Cobalt, total recoverable | <6.21 | JU | V | < EQL | <20 | U | | < EQL | ug/L |
| | Copper, total recoverable | <7.16 | U | V | < EQL | <3.76 | JU | | < EQL | ug/L |
| | Cyanide | <10 | U | | < EQL | | | | | ug/L |
| | Di-n-butyl phthalate | <11 | U | | < EQL | <9.8 | U | | < EQL | ug/L |
| | Di-n-octyl phthalate | <11 | U | | < EQL | | | | | ug/L |
| | Diallate | <11 | U | | < EQL | | | | | ug/L |
| | Dibenz(a,h)anthracene | <11 | U | | < EQL | | | | | ug/L |
| | Dibenzofuran | <11 | U | | < EQL | | | | | ug/L |
| | Dibromochloromethane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Dibromomethane (Methylene bromide) | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Dichlorodifluoromethane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Dichloromethane (Methylene chloride) | <2.7 | U | V | < EQL | <10 | U | | < EQL | ug/L |
| | Dieldrin | <.21 | U | | < EQL | | | | | ug/L |
| | Diethyl phthalate | <11 | U | | < EQL | | | | | ug/L |
| | Dimethoate | <11 | U | | < EQL | | | | | ug/L |
| | Dimethyl phthalate | <11 | U | | < EQL | | | | | ug/L |
| | Diphenylamine | <11 | U | | < EQL | | | | | ug/L |
| | Disulfoton | <11 | U | | < EQL | | | | | ug/L |
| | Endosulfan I | <.1 | U | | < EQL | | | | | ug/L |
| | Endosulfan II | <.21 | U | | < EQL | | | | | ug/L |
| | Endosulfan sulfate | <.21 | U | | < EQL | | | | | ug/L |
| | Endrin | <.21 | U | | < EQL | | | | | ug/L |

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+ = exceeded the Groundwater Protection Standards listed in Appendix A.

Table B. Assessment Monitoring Results for Groundwater Wells (Cont.)

WELL LFW 76

ANALYTICAL DATA

Groundwater Protection Standard

| ST | Parameter | 1st Half | CLP | EPA | Filt. | 2nd Half | CLP | EPA | Filt. | Unit |
|----|---------------------------------|----------|-----|-----|-------|----------|-----|-----|-------|-------|
| | Endrin aldehyde | <.21 | U | | < EQL | | | | | ug/L |
| | Ethyl methacrylate | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Ethyl methanesulfonate | <11 | U | | < EQL | | | | | ug/L |
| | Ethylbenzene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Famphur | <200 | U | | < EQL | | | | | ug/L |
| | Fluoranthene | <11 | U | | < EQL | | | | | ug/L |
| | Fluorene | <11 | U | | < EQL | | | | | ug/L |
| | Heptachlor | <.1 | U | | < EQL | | | | | ug/L |
| | Heptachlor epoxide | <.1 | U | | < EQL | | | | | ug/L |
| | Hexachlorobenzene | <11 | U | | < EQL | | | | | ug/L |
| | Hexachlorobutadiene | <11 | U | | < EQL | | | | | ug/L |
| | Hexachlorocyclopentadiene | <11 | U | | < EQL | | | | | ug/L |
| | Hexachloroethane | <11 | U | | < EQL | | | | | ug/L |
| | Hexachloropropene | <52 | U | | < EQL | | | | | ug/L |
| | Indeno(1,2,3-c,d)pyrene | <11 | U | | < EQL | | | | | ug/L |
| | Iodomethane (Methyl iodide) | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Isobutyl alcohol | <500 | U | | < EQL | <500 | U | | < EQL | ug/L |
| | Isodrin | <11 | U | | < EQL | | | | | ug/L |
| | Isophorone | <11 | U | | < EQL | | | | | ug/L |
| | Isosafrole | <11 | U | | < EQL | | | | | ug/L |
| | Kepone | <11 | U | | < EQL | | | | | ug/L |
| | Lead, total recoverable | 3.51 | J | I | NDD | <10 | U | | < EQL | ug/L |
| | Lindane | <.1 | U | | < EQL | | | | | ug/L |
| | Mercury, total recoverable | <.5 | U | | < EQL | <.5 | U | | < EQL | ug/L |
| | Methacrylonitrile | <200 | U | | < EQL | <200 | U | | < EQL | ug/L |
| | Methapyrilene | <11 | U | | < EQL | | | | | ug/L |
| | Methoxychlor | <.1 | U | | < EQL | | | | | ug/L |
| | Methyl ethyl ketone | <20 | U | | < EQL | <20 | U | | < EQL | ug/L |
| | Methyl isobutyl ketone | <10 | U | | < EQL | <10 | U | | < EQL | ug/L |
| | Methyl methacrylate | <20 | U | | < EQL | <20 | U | | < EQL | ug/L |
| | Methyl methanesulfonate | <11 | U | | < EQL | | | | | ug/L |
| | N-Nitrosodi-n-butylamine | <11 | U | | < EQL | | | | | ug/L |
| | N-Nitrosodiethylamine | <11 | U | | < EQL | | | | | ug/L |
| | N-Nitrosodimethylamine | <26 | U | | < EQL | | | | | ug/L |
| | N-Nitrosodiphenylamine | <11 | U | | < EQL | | | | | ug/L |
| | N-Nitrosodipropylamine | <11 | U | | < EQL | | | | | ug/L |
| | N-Nitrosomethylethylamine | <11 | U | | < EQL | | | | | ug/L |
| | N-Nitrosopiperidine | <11 | U | | < EQL | | | | | ug/L |
| | N-Nitrosopyrrolidine | <11 | U | | < EQL | | | | | ug/L |
| | Naphthalene | <11 | U | | < EQL | | | | | ug/L |
| | Nickel, total recoverable | <10 | JU | | < EQL | <10.9 | JU | | < EQL | ug/L |
| | Nitrobenzene | <11 | U | | < EQL | | | | | ug/L |
| | O,O,O-Triethyl phosphorothioate | <11 | U | | < EQL | | | | | ug/L |
| | PCB | | | | | | | | | |
| | PCB 1016 | <1 | U | | < EQL | | | | | ug/L |
| | PCB 1232 | <1 | U | | < EQL | | | | | ug/L |
| | PCB 1242 | <2.1 | U | | < EQL | | | | | ug/L |
| | PCB 1248 | <1 | U | | < EQL | | | | | ug/L |
| | PCB 1254 | <1 | U | | < EQL | | | | | ug/L |
| | PCB 1260 | <1 | U | | < EQL | | | | | ug/L |
| | Parathion | <11 | U | | < EQL | | | | | ug/L |
| | Parathion methyl | <11 | U | | < EQL | | | | | ug/L |
| | Pentachlorobenzene | <52 | U | | < EQL | | | | | ug/L |
| | Pentachloronitrobenzene | <11 | U | | < EQL | | | | | ug/L |
| | Pentachlorophenol | <26 | U | | < EQL | | | | | ug/L |
| | Phenacetin | <11 | U | | < EQL | | | | | ug/L |
| | Phenanthrene | <11 | U | | < EQL | | | | | ug/L |
| | Phenol | <11 | U | | < EQL | | | | | ug/L |
| | Phorate | <11 | U | | < EQL | | | | | ug/L |
| | Pronamid | <11 | U | | < EQL | | | | | ug/L |
| | Propionitrile | <200 | U | | < EQL | <200 | U | | < EQL | ug/L |
| | Pyrene | <11 | U | | < EQL | | | | | ug/L |
| | Safrole | <11 | U | | < EQL | | | | | ug/L |
| | Selenium, total recoverable | <10 | U | | < EQL | <10 | U | | < EQL | ug/L |
| | Silver, total recoverable | <4.74 | JU | | < EQL | <20 | U | | < EQL | ug/L |
| | Specific conductance | 24.2 | | | < 37 | 20.6 | | | < 37 | uS/cm |
| | Styrene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Sulfide | <1000 | U | | < EQL | | | | | ug/L |

Notes: Concentrations in bold italics exceed the groundwater protection standards listed in Appendix A. Bold italics were not assigned if the result was qualified with a "J", "R", "L", or "U" qualifier.

+ = exceeded the Groundwater Protection Standards listed in Appendix A.

Table B. Assessment Monitoring Results for Groundwater Wells (Cont.)

WELL LFW 76

ANALYTICAL DATA

Groundwater Protection Standard

| ST | Parameter | 1st Half | CLP | EPA | Filt. | 2nd Half | CLP | EPA | Filt. | Unit |
|----|-----------------------------|-------------|-----|-----|--------|----------|-----|-----|-------|------|
| | Tetrachloroethylene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Thallium, total recoverable | <3.57 | JU | | < EQL | <10 | U | | < EQL | ug/L |
| | Thionazin | <11 | U | | < EQL | | | | | ug/L |
| | Tin, total recoverable | <200 | U | | < EQL | | | | | ug/L |
| | Toluene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Toxaphene | <2.1 | U | | < EQL | | | | | ug/L |
| | Trichloroethylene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Trichlorofluoromethane | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Vanadium, total recoverable | <2.36 | JU | | < EQL | 7.24 | J | I | NDD | ug/L |
| | Vinyl acetate | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | Xylenes | <10 | U | | < EQL | <10 | U | | < EQL | ug/L |
| | Zinc, total recoverable | 35.6 | | | > 29.3 | <9.95 | JU | | < EQL | ug/L |
| | alpha-Benzene hexachloride | <.1 | U | | < EQL | | | | | ug/L |
| | alpha-Chlordane | <.1 | U | | < EQL | | | | | ug/L |
| | beta-Benzene hexachloride | <.1 | U | | < EQL | <.1 | U | | < EQL | ug/L |
| | cis-1,2-Dichloroethylene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | cis-1,3-Dichloropropene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | delta-Benzene hexachloride | <.1 | U | | < EQL | | | | | ug/L |
| | gamma-Chlordane | <.1 | U | | < EQL | | | | | ug/L |
| | m-Cresol (3-Methylphenol) | | | | | | | | | |
| | m-Nitroaniline | <26 | U | | < EQL | | | | | ug/L |
| | o-Cresol (2-Methylphenol) | <11 | U | | < EQL | | | | | ug/L |
| | o-Nitroaniline | <26 | U | | < EQL | | | | | ug/L |
| | o-Toluidine | <11 | U | | < EQL | | | | | ug/L |
| | p,p"-DDD | <.21 | U | | < EQL | | | | | ug/L |
| | p,p"-DDE | <.21 | U | | < EQL | | | | | ug/L |
| | p,p"-DDT | <.21 | U | | < EQL | | | | | ug/L |
| | p-Cresol (4-Methylphenol) | <10 | U | | < EQL | | | | | ug/L |
| | p-Dimethylaminoazobenzene | <11 | U | | < EQL | | | | | ug/L |
| | p-Nitroaniline | <11 | U | | < EQL | | | | | ug/L |
| | p-Phenylenediamine | <11 | U | | < EQL | | | | | ug/L |
| | pH | 5.63 | J | Q | NDD | 5.14 | J | Q | NDD | pH |
| | trans-1,2-Dichloroethylene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | trans-1,3-Dichloropropene | <5 | U | | < EQL | <5 | U | | < EQL | ug/L |
| | trans-1,4-Dichloro-2-butene | <20 | U | | < EQL | <20 | U | | < EQL | ug/L |

Notes: Concentrations in bold italics exceed the groundwater protection standards listed in Appendix A. Bold italics were not assigned if the result was qualified with a "J", "R", "L", or "U" qualifier.

+ = exceeded the Groundwater Protection Standards listed in Appendix A.

Table B. Assessment Monitoring Results for Groundwater Wells (Cont.)

WELL LFW 78

| <u>SRS Coord.</u> | <u>Lat/Longitude</u> | <u>Screen Zone Elevation</u> | <u>Top of Casing</u> | <u>Casing</u> | <u>Pump</u> | <u>Screen Zone</u> |
|------------------------|----------------------------------|------------------------------|----------------------|---------------|-------------|--------------------|
| N 86064.9 E 44726.5 | 33.28909 Deg N -81.7162 Deg W | 164.92 - 149.9 ft msl | 238.92 ft msl | 2 " PVC | V | Upper |

SAMPLE DATE 03/30/00 09/29/00

FIELD DATA

| <u>Parameter</u> | <u>1st Half</u> | <u>2nd Half</u> | <u>Unit</u> |
|---------------------|-----------------|-----------------|-------------|
| Water Elevation | 159.62 | 158.66 | ft msl |
| pH | 5 | 4.8 | pH |
| Sp. Conductance | 38 | 35 | uS/cm |
| Water temperature | 26.7 | 23.7 | deg. C |
| Alkalinity as CaCO3 | 5 | 0 | mg/L |
| Turbidity | 9.8 | 8.5 | NTU |
| Volumes purged | 6.39 | 4.26 | well volume |
| Sampling code | | NX | |

ANALYTICAL DATA

Groundwater Protection Standard

| <u>SI</u> | <u>Parameter</u> | <u>1st Half</u> | <u>CLP</u> | <u>EPA</u> | <u>Filt.</u> | <u>2nd Half</u> | <u>CLP</u> | <u>EPA</u> | <u>Filt.</u> | <u>Unit</u> |
|-----------|--------------------------------|-----------------|------------|------------|--------------|-----------------|------------|------------|--------------|-------------|
| | 1,1,1,2-Tetrachloroethane | <5 | JU | L | < EQL | <5 | U | | < EQL | ug/L |
| | 1,1,1-Trichloroethane | <5 | JU | L | < EQL | <5 | U | | < EQL | ug/L |
| | 1,1,2,2-Tetrachloroethane | <5 | JU | L | < EQL | <5 | U | | < EQL | ug/L |
| | 1,1,2-Trichloroethane | <5 | JU | L | < EQL | <5 | U | | < EQL | ug/L |
| | 1,1-Dichloroethane | <5 | JU | L | < EQL | <5 | U | | < EQL | ug/L |
| | 1,1-Dichloroethylene | <5 | JU | L | < EQL | <5 | U | | < EQL | ug/L |
| | 1,1-Dichloropropene | <5 | JU | L | < EQL | <5 | U | | < EQL | ug/L |
| | 1,2,3-Trichloropropane | <5 | JU | L | < EQL | <5 | U | | < EQL | ug/L |
| | 1,2,4,5-Tetrachlorobenzene | <52 | JU | Q | < EQL | | | | | ug/L |
| | 1,2,4-Trichlorobenzene | <10 | JU | Q | < EQL | | | | | ug/L |
| | 1,2-Dibromo-3-chloropropane | <10 | JU | L | < EQL | <10 | U | | < EQL | ug/L |
| | 1,2-Dibromoethane | <5 | JU | L | < EQL | <5 | U | | < EQL | ug/L |
| | 1,2-Dichlorobenzene | <10 | JU | Q | < EQL | <5 | U | | < EQL | ug/L |
| | 1,2-Dichloroethane | <5 | JU | L | < EQL | <5 | U | | < EQL | ug/L |
| | 1,2-Dichloropropane | <5 | JU | L | < EQL | <5 | U | | < EQL | ug/L |
| | 1,3,5-Trinitrobenzene | <10 | JU | Q | < EQL | | | | | ug/L |
| | 1,3-Dichlorobenzene | <10 | JU | Q | < EQL | <5 | U | | < EQL | ug/L |
| | 1,3-Dichloropropane | <5 | JU | L | < EQL | <5 | U | | < EQL | ug/L |
| | 1,3-Dinitrobenzene | <50 | JU | Q | < EQL | | | | | ug/L |
| | 1,4-Dichlorobenzene | <10 | JU | Q | < EQL | <5 | U | | < EQL | ug/L |
| | 1,4-Naphthoquinone | <10 | JU | Q | < EQL | | | | | ug/L |
| | 1-Naphthylamine | <10 | JU | Q | < EQL | | | | | ug/L |
| | 2,2-Dichloropropane | <5 | JU | L | < EQL | <5 | U | | < EQL | ug/L |
| | 2,2-Oxybis(1-chloropropane) | <10 | JU | Q | < EQL | | | | | ug/L |
| | 2,3,4,6-Tetrachlorophenol | <10 | JU | Q | < EQL | | | | | ug/L |
| | 2,4,5-T | <2 | JU | Q | < EQL | | | | | ug/L |
| | 2,4,5-TP (Silvex) | <2 | JU | Q | < EQL | | | | | ug/L |
| | 2,4,5-Trichlorophenol | <10 | JU | Q | < EQL | | | | | ug/L |
| | 2,4,6-Trichlorophenol | <26 | JU | Q | < EQL | | | | | ug/L |
| | 2,4-Dichlorophenol | <10 | JU | Q | < EQL | | | | | ug/L |
| | 2,4-Dichlorophenoxyacetic acid | <2 | JU | Q | < EQL | | | | | ug/L |
| | 2,4-Dimethyl phenol | <10 | JU | Q | < EQL | | | | | ug/L |
| | 2,4-Dinitrophenol | <26 | JU | Q | < EQL | | | | | ug/L |
| | 2,4-Dinitrotoluene | <10 | JU | Q | < EQL | | | | | ug/L |
| | 2,6-Dichlorophenol | <25 | JU | Q | < EQL | | | | | ug/L |
| | 2,6-Dinitrotoluene | <10 | JU | Q | < EQL | | | | | ug/L |
| | 2-Acetylaminofluorene | <10 | JU | Q | < EQL | | | | | ug/L |
| | 2-Chloronaphthalene | <10 | JU | Q | < EQL | | | | | ug/L |
| | 2-Chlorophenol | <10 | JU | Q | < EQL | | | | | ug/L |
| | 2-Hexanone | <20 | JU | L | < EQL | <20 | U | | < EQL | ug/L |
| | 2-Methyl-4,6-dinitrophenol | <26 | JU | Q | < EQL | | | | | ug/L |
| | 2-Methylnaphthalene | <10 | JU | Q | < EQL | | | | | ug/L |
| | 2-Naphthylamine | <10 | JU | Q | < EQL | | | | | ug/L |
| | 2-Nitrophenol | <10 | JU | Q | < EQL | | | | | ug/L |
| | 2-sec-Butyl-4,6-dinitrophenol | <2 | JU | LQ | < EQL | | | | | ug/L |
| | 3,3'-Dichlorobenzidine | <10 | JU | Q | < EQL | | | | | ug/L |
| | 3,3'-Dimethylbenzidine | <21 | JU | Q | < EQL | | | | | ug/L |
| | 3-Methylcholanthrene | <10 | JU | Q | < EQL | | | | | ug/L |
| | 4-Aminobiphenyl | <10 | JU | Q | < EQL | | | | | ug/L |
| | 4-Bromophenyl phenyl ether | <10 | JU | Q | < EQL | | | | | ug/L |
| | 4-Chloro-m-cresol | <10 | JU | Q | < EQL | | | | | ug/L |

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+ = exceeded the Groundwater Protection Standards listed in Appendix A.

Table B. Assessment Monitoring Results for Groundwater Wells (Cont.)

WELL LFW 78

ANALYTICAL DATA

Groundwater Protection Standard

| ST | Parameter | 1st Half | CLP | EPA | Filt. | 2nd Half | CLP | EPA | Filt. | Unit |
|----|--------------------------------------|----------|-----|-----|-------|----------|-----|-----|--------|------|
| | 4-Chloroaniline | <10 | JU | Q | < EQL | | | | | ug/L |
| | 4-Chlorophenyl phenyl ether | <10 | JU | Q | < EQL | | | | | ug/L |
| | 4-Nitrophenol | <26 | JU | Q | < EQL | | | | | ug/L |
| | 5-Nitro-o-toluidine | <10 | JU | Q | < EQL | | | | | ug/L |
| | 7,12-Dimethylbenz(a)anthracene | <10 | JU | Q | < EQL | | | | | ug/L |
| | Acenaphthene | <10 | JU | Q | < EQL | | | | | ug/L |
| | Acenaphthylene | <10 | JU | Q | < EQL | | | | | ug/L |
| | Acetone | <20 | JU | L | < EQL | <20 | U | | < EQL | ug/L |
| | Acetonitrile (Methyl cyanide) | <200 | JU | L | < EQL | <200 | U | | < EQL | ug/L |
| | Acetophenone | <10 | JU | Q | < EQL | | | | | ug/L |
| | Acrolein | <50 | JU | L | < EQL | <50 | U | | < EQL | ug/L |
| | Acrylonitrile | <10 | JU | L | < EQL | <10 | U | | < EQL | ug/L |
| | Aldrin | <.11 | JU | Q | < EQL | | | | | ug/L |
| | Allyl chloride | <5 | JU | L | < EQL | <5 | U | | < EQL | ug/L |
| | Anthracene | <10 | JU | Q | < EQL | | | | | ug/L |
| | Antimony, total recoverable | <100 | U | | < EQL | <100 | U | | < EQL | ug/L |
| | Arsenic, total recoverable | <10 | U | | < EQL | <10 | U | | < EQL | ug/L |
| | Barium, total recoverable | 9.43 | J | I | NDD | 9.17 | J | I | NDD | ug/L |
| | Benzene | <5 | JU | L | < EQL | <5 | U | | < EQL | ug/L |
| | Benzo(a)anthracene | <10 | JU | Q | < EQL | | | | | ug/L |
| | Benzo(a)pyrene | <10 | JU | Q | < EQL | | | | | ug/L |
| | Benzo(b)fluoranthene | <10 | JU | Q | < EQL | | | | | ug/L |
| | Benzo(g,h,i)perylene | <10 | JU | Q | < EQL | | | | | ug/L |
| | Benzo(k)fluoranthene | <10 | JU | Q | < EQL | | | | | ug/L |
| | Benzyl alcohol | | | | | | | | | |
| | Beryllium, total recoverable | <1 | U | | < EQL | <.12 | JU | | < EQL | ug/L |
| | Bis(2-chloroethoxy) methane | <10 | JU | Q | < EQL | | | | | ug/L |
| | Bis(2-chloroethyl) ether | <10 | JU | Q | < EQL | | | | | ug/L |
| | Bis(2-ethylhexyl) phthalate | <3.9 | JU | IQV | < EQL | | | | | ug/L |
| | Bromochloromethane | <5 | JU | L | < EQL | <5 | U | | < EQL | ug/L |
| | Bromodichloromethane | <5 | JU | L | < EQL | <5 | U | | < EQL | ug/L |
| | Bromoform | <5 | JU | L | < EQL | <5 | U | | < EQL | ug/L |
| | Bromomethane (Methyl bromide) | <5 | JU | L | < EQL | <5 | U | | < EQL | ug/L |
| | Butylbenzyl phthalate | <10 | JU | Q | < EQL | | | | | ug/L |
| | Cadmium, total recoverable | <10 | U | | < EQL | <10 | U | | < EQL | ug/L |
| | Carbon disulfide | <5 | JU | L | < EQL | <5 | U | | < EQL | ug/L |
| | Carbon tetrachloride | <5 | JU | L | < EQL | <5 | U | | < EQL | ug/L |
| | Chlorobenzene | <5 | JU | L | < EQL | <5 | U | | < EQL | ug/L |
| | Chlorobenzilate | <10 | JU | Q | < EQL | | | | | ug/L |
| | Chloroethane | <5 | JU | L | < EQL | <5 | U | | < EQL | ug/L |
| | Chloroethene (Vinyl chloride) | <5 | JU | L | < EQL | <5 | U | | < EQL | ug/L |
| | Chloroform | <5 | JU | L | < EQL | <5 | U | | < EQL | ug/L |
| | Chloromethane (Methyl chloride) | <5 | JU | L | < EQL | <5 | U | | < EQL | ug/L |
| | Chloroprene | <20 | JU | L | < EQL | <20 | U | | < EQL | ug/L |
| | Chromium, total recoverable | <10 | U | | < EQL | 10.1 | | | < 100 | ug/L |
| | Chrysene | <10 | JU | Q | < EQL | | | | | ug/L |
| | Cobalt, total recoverable | <20 | U | | < EQL | <20 | U | | < EQL | ug/L |
| | Copper, total recoverable | 9.42 | J | I | NDD | 24.5 | | | < 29.8 | ug/L |
| | Cyanide | <10 | JU | Q | < EQL | | | | | ug/L |
| | Di-n-butyl phthalate | <10 | JU | Q | < EQL | <11 | JU | Q | < EQL | ug/L |
| | Di-n-octyl phthalate | <10 | JU | Q | < EQL | | | | | ug/L |
| | Diallate | <10 | JU | Q | < EQL | | | | | ug/L |
| | Dibenz(a,h)anthracene | <10 | JU | Q | < EQL | | | | | ug/L |
| | Dibenzofuran | <10 | JU | Q | < EQL | | | | | ug/L |
| | Dibromochloromethane | <5 | JU | L | < EQL | <5 | U | | < EQL | ug/L |
| | Dibromomethane (Methylene bromide) | <5 | JU | L | < EQL | <5 | U | | < EQL | ug/L |
| | Dichlorodifluoromethane | <5 | JU | L | < EQL | <5 | U | | < EQL | ug/L |
| | Dichloromethane (Methylene chloride) | <10 | JU | L | < EQL | <10 | U | | < EQL | ug/L |
| | Dieldrin | <.21 | JU | Q | < EQL | | | | | ug/L |
| | Diethyl phthalate | <10 | JU | Q | < EQL | | | | | ug/L |
| | Dimethoate | <10 | JU | Q | < EQL | | | | | ug/L |
| | Dimethyl phthalate | <10 | JU | Q | < EQL | | | | | ug/L |
| | Diphenylamine | <10 | JU | Q | < EQL | | | | | ug/L |
| | Disulfoton | <10 | JU | Q | < EQL | | | | | ug/L |
| | Endosulfan I | <.11 | JU | Q | < EQL | | | | | ug/L |
| | Endosulfan II | <.21 | JU | Q | < EQL | | | | | ug/L |
| | Endosulfan sulfate | <.21 | JU | Q | < EQL | | | | | ug/L |
| | Endrin | <.21 | JU | Q | < EQL | | | | | ug/L |

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+ = exceeded the Groundwater Protection Standards listed in Appendix A.

Table B. Assessment Monitoring Results for Groundwater Wells (Cont.)

WELL LFW 78

ANALYTICAL DATA

Groundwater Protection Standard

| ST | Parameter | 1st Half | CLP | EPA | Filt. | 2nd Half | CLP | EPA | Filt. | Unit |
|----|---------------------------------|-----------|-----|-----|-------|----------|-----|-----|-------|-------|
| | Endrin aldehyde | <.21 | JU | Q | < EQL | | | | | ug/L |
| | Ethyl methacrylate | <5 | JU | L | < EQL | <5 | U | | < EQL | ug/L |
| | Ethyl methanesulfonate | <10 | JU | Q | < EQL | | | | | ug/L |
| | Ethylbenzene | <5 | JU | L | < EQL | <5 | U | | < EQL | ug/L |
| | Famphur | <200 | JU | Q | < EQL | | | | | ug/L |
| | Fluoranthene | <10 | JU | Q | < EQL | | | | | ug/L |
| | Fluorene | <10 | JU | Q | < EQL | | | | | ug/L |
| | Heptachlor | <.11 | JU | Q | < EQL | | | | | ug/L |
| | Heptachlor epoxide | <.11 | JU | Q | < EQL | | | | | ug/L |
| | Hexachlorobenzene | <10 | JU | Q | < EQL | | | | | ug/L |
| | Hexachlorobutadiene | <10 | JU | Q | < EQL | | | | | ug/L |
| | Hexachlorocyclopentadiene | <10 | JU | Q | < EQL | | | | | ug/L |
| | Hexachloroethane | <10 | JU | Q | < EQL | | | | | ug/L |
| | Hexachloropropene | <52 | JU | Q | < EQL | | | | | ug/L |
| | Indeno(1,2,3-c,d)pyrene | <10 | JU | Q | < EQL | | | | | ug/L |
| | Iodomethane (Methyl iodide) | <5 | JU | L | < EQL | <5 | U | | < EQL | ug/L |
| | Isobutyl alcohol | <500 | JU | L | < EQL | <500 | U | | < EQL | ug/L |
| | Isodrin | <10 | JU | Q | < EQL | | | | | ug/L |
| | Isophorone | <10 | JU | Q | < EQL | | | | | ug/L |
| | Isosafrole | <10 | JU | Q | < EQL | | | | | ug/L |
| | Kepone | <10 | JU | Q | < EQL | | | | | ug/L |
| | Lead, total recoverable | <10 | U | | < EQL | <10 | U | | < EQL | ug/L |
| | Lindane | <.11 | JU | Q | < EQL | | | | | ug/L |
| | Mercury, total recoverable | <.5 | U | | < EQL | <.5 | U | | < EQL | ug/L |
| | Methacrylonitrile | <200 | JU | L | < EQL | <200 | U | | < EQL | ug/L |
| | Methapyrene | <10 | JU | Q | < EQL | | | | | ug/L |
| | Methoxychlor | <1.1 | JU | Q | < EQL | | | | | ug/L |
| | Methyl ethyl ketone | <20 | JU | L | < EQL | <20 | U | | < EQL | ug/L |
| | Methyl isobutyl ketone | <10 | JU | L | < EQL | <10 | U | | < EQL | ug/L |
| | Methyl methacrylate | <20 | JU | L | < EQL | <20 | U | | < EQL | ug/L |
| | Methyl methanesulfonate | <10 | JU | Q | < EQL | | | | | ug/L |
| | N-Nitrosodi-n-butylamine | <10 | JU | Q | < EQL | | | | | ug/L |
| | N-Nitrosodiethylamine | <10 | JU | Q | < EQL | | | | | ug/L |
| | N-Nitrosodimethylamine | <26 | JU | Q | < EQL | | | | | ug/L |
| | N-Nitrosodiphenylamine | <10 | JU | Q | < EQL | | | | | ug/L |
| | N-Nitrosodipropylamine | <10 | JU | Q | < EQL | | | | | ug/L |
| | N-Nitrosomethylethylamine | <10 | JU | Q | < EQL | | | | | ug/L |
| | N-Nitrosopiperidine | <10 | JU | Q | < EQL | | | | | ug/L |
| | N-Nitrosopyrrolidine | <10 | JU | Q | < EQL | | | | | ug/L |
| | Naphthalene | <10 | JU | Q | < EQL | | | | | ug/L |
| | Nickel, total recoverable | <12.7 | JU | I | < EQL | 16.6 | J | I | NDD | ug/L |
| | Nitrobenzene | <10 | JU | Q | < EQL | | | | | ug/L |
| | O,O,O-Triethyl phosphorothioate | <10 | JU | Q | < EQL | | | | | ug/L |
| | PCB | | | | | | | | | |
| | PCB 1016 | <1.1 | JU | Q | < EQL | | | | | ug/L |
| | PCB 1232 | <1.1 | JU | Q | < EQL | | | | | ug/L |
| | PCB 1242 | <2.1 | JU | Q | < EQL | | | | | ug/L |
| | PCB 1248 | <1.1 | JU | Q | < EQL | | | | | ug/L |
| | PCB 1254 | <1.1 | JU | Q | < EQL | | | | | ug/L |
| | PCB 1260 | <1.1 | JU | Q | < EQL | | | | | ug/L |
| | Parathion | <10 | JU | Q | < EQL | | | | | ug/L |
| | Parathion methyl | <10 | JU | Q | < EQL | | | | | ug/L |
| | Pentachlorobenzene | <52 | JU | Q | < EQL | | | | | ug/L |
| | Pentachloronitrobenzene | <10 | JU | Q | < EQL | | | | | ug/L |
| | Pentachlorophenol | <26 | JU | Q | < EQL | | | | | ug/L |
| | Phenacetin | <10 | JU | Q | < EQL | | | | | ug/L |
| | Phenanthrene | <10 | JU | Q | < EQL | | | | | ug/L |
| | Phenol | <10 | JU | Q | < EQL | | | | | ug/L |
| | Phorate | <10 | JU | Q | < EQL | | | | | ug/L |
| | Pronamid | <10 | JU | Q | < EQL | | | | | ug/L |
| | Propionitrile | <200 | JU | L | < EQL | <200 | U | | < EQL | ug/L |
| | Pyrene | <10 | JU | Q | < EQL | | | | | ug/L |
| | Safrole | <10 | JU | Q | < EQL | | | | | ug/L |
| | Selenium, total recoverable | <10 | U | | < EQL | <10 | U | | < EQL | ug/L |
| | Silver, total recoverable | <20 | U | | < EQL | <3.49 | JU | | < EQL | ug/L |
| | Specific conductance | 39 | | | > 37 | 33 | | | < 37 | uS/cm |
| | Styrene | <5 | JU | L | < EQL | <5 | U | | < EQL | ug/L |
| | Sulfide | <1000 | U | | < EQL | | | | | ug/L |

Notes: Concentrations in bold italics exceed the groundwater protection standards listed in Appendix A. Bold Italics were not assigned if the result was qualified with a "J", "R", "L", or "U" qualifier.

+ = exceeded the Groundwater Protection Standards listed in Appendix A.

Table B. Assessment Monitoring Results for Groundwater Wells (Cont.)

WELL LFW 78

ANALYTICAL DATA

Groundwater Protection Standard

| ST | Parameter | 1st Half | CLP | EPA | Filt. | 2nd Half | CLP | EPA | Filt. | Unit |
|----|-----------------------------|-------------|-----|-----|--------|-------------|-----|-----|--------|------|
| | Tetrachloroethylene | <5 | JU | L | < EQL | <5 | U | | < EQL | ug/L |
| | Thallium, total recoverable | <10 | U | | < EQL | <10 | U | | < EQL | ug/L |
| | Thionazin | <10 | JU | Q | < EQL | | | | | ug/L |
| | Tin, total recoverable | <200 | U | | < EQL | | | | | ug/L |
| | Toluene | <5 | JU | L | < EQL | <5 | U | | < EQL | ug/L |
| | Toxaphene | <2.1 | JU | Q | < EQL | | | | | ug/L |
| | Trichloroethylene | <5 | JU | L | < EQL | <5 | U | | < EQL | ug/L |
| | Trichlorofluoromethane | 27 | J | L | NDD | 25 | J | K | NDD | ug/L |
| | Vanadium, total recoverable | <10 | U | | < EQL | <10 | U | | < EQL | ug/L |
| | Vinyl acetate | <5 | JU | L | < EQL | <5 | U | | < EQL | ug/L |
| | Xylenes | <10 | JU | L | < EQL | <10 | U | | < EQL | ug/L |
| + | Zinc, total recoverable | 32.7 | | | > 29.3 | 29.4 | | | > 29.3 | ug/L |
| | alpha-Benzene hexachloride | <.11 | JU | Q | < EQL | | | | | ug/L |
| | alpha-Chlordane | <.11 | JU | Q | < EQL | | | | | ug/L |
| | beta-Benzene hexachloride | <.11 | JU | Q | < EQL | <.1 | U | | < EQL | ug/L |
| | cis-1,2-Dichloroethylene | <5 | JU | L | < EQL | <5 | U | | < EQL | ug/L |
| | cis-1,3-Dichloropropene | <5 | JU | L | < EQL | <5 | U | | < EQL | ug/L |
| | delta-Benzene hexachloride | <.11 | JU | Q | < EQL | | | | | ug/L |
| | gamma-Chlordane | <.11 | JU | Q | < EQL | | | | | ug/L |
| | m-Cresol (3-Methylphenol) | | | | | | | | | |
| | m-Nitroaniline | <26 | JU | Q | < EQL | | | | | ug/L |
| | o-Cresol (2-Methylphenol) | <10 | JU | Q | < EQL | | | | | ug/L |
| | o-Nitroaniline | <26 | JU | Q | < EQL | | | | | ug/L |
| | o-Toluidine | <10 | JU | Q | < EQL | | | | | ug/L |
| | p,p"-DDD | <.21 | JU | Q | < EQL | | | | | ug/L |
| | p,p"-DDE | <.21 | JU | Q | < EQL | | | | | ug/L |
| | p,p"-DDT | <.21 | JU | Q | < EQL | | | | | ug/L |
| | p-Cresol (4-Methylphenol) | <10 | JU | Q | < EQL | | | | | ug/L |
| | p-Dimethylaminoazobenzene | <10 | JU | Q | < EQL | | | | | ug/L |
| | p-Nitroaniline | <10 | JU | Q | < EQL | | | | | ug/L |
| | p-Phenylenediamine | <10 | JU | Q | < EQL | | | | | ug/L |
| | pH | 5.3 | J | Q | NDD | 4.93 | J | Q | NDD | pH |
| | trans-1,2-Dichloroethylene | <5 | JU | L | < EQL | <5 | U | * | < EQL | ug/L |
| | trans-1,3-Dichloropropene | <5 | JU | L | < EQL | <5 | U | | < EQL | ug/L |
| | trans-1,4-Dichloro-2-butene | <20 | JU | L | < EQL | <20 | U | | < EQL | ug/L |

Notes: Concentrations in bold italics exceed the groundwater protection standards listed in Appendix A. Bold italics were not assigned if the result was qualified with a "J", "R", "L", or "U" qualifier.

+ = exceeded the Groundwater Protection Standards listed in Appendix A.

Appendix C

Figures

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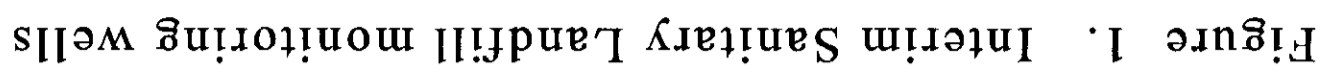
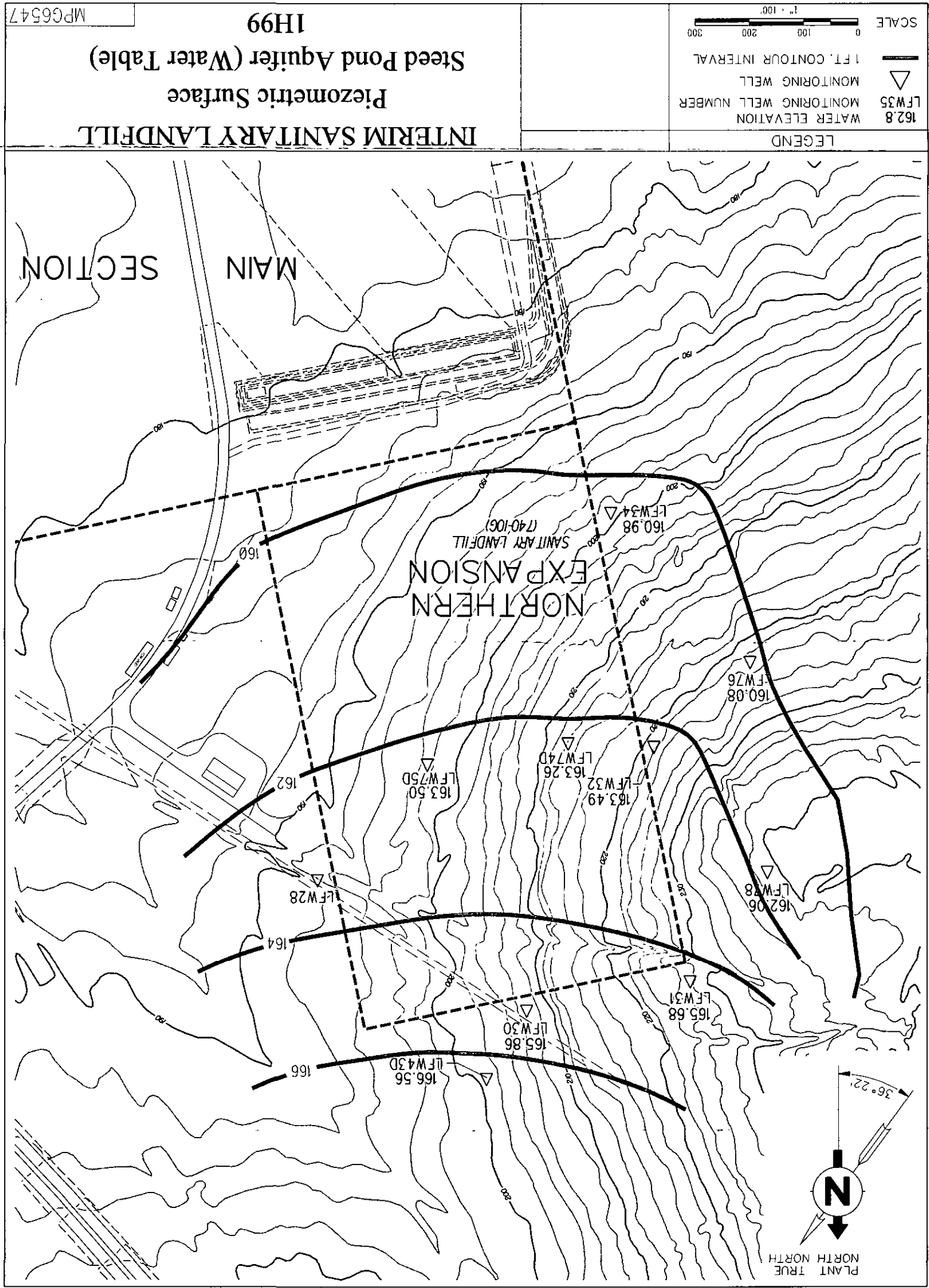


Figure 2. Potentiometric surface map of the Steed Pond Aquifer, 1H99



86826-44434

86826 -
44434

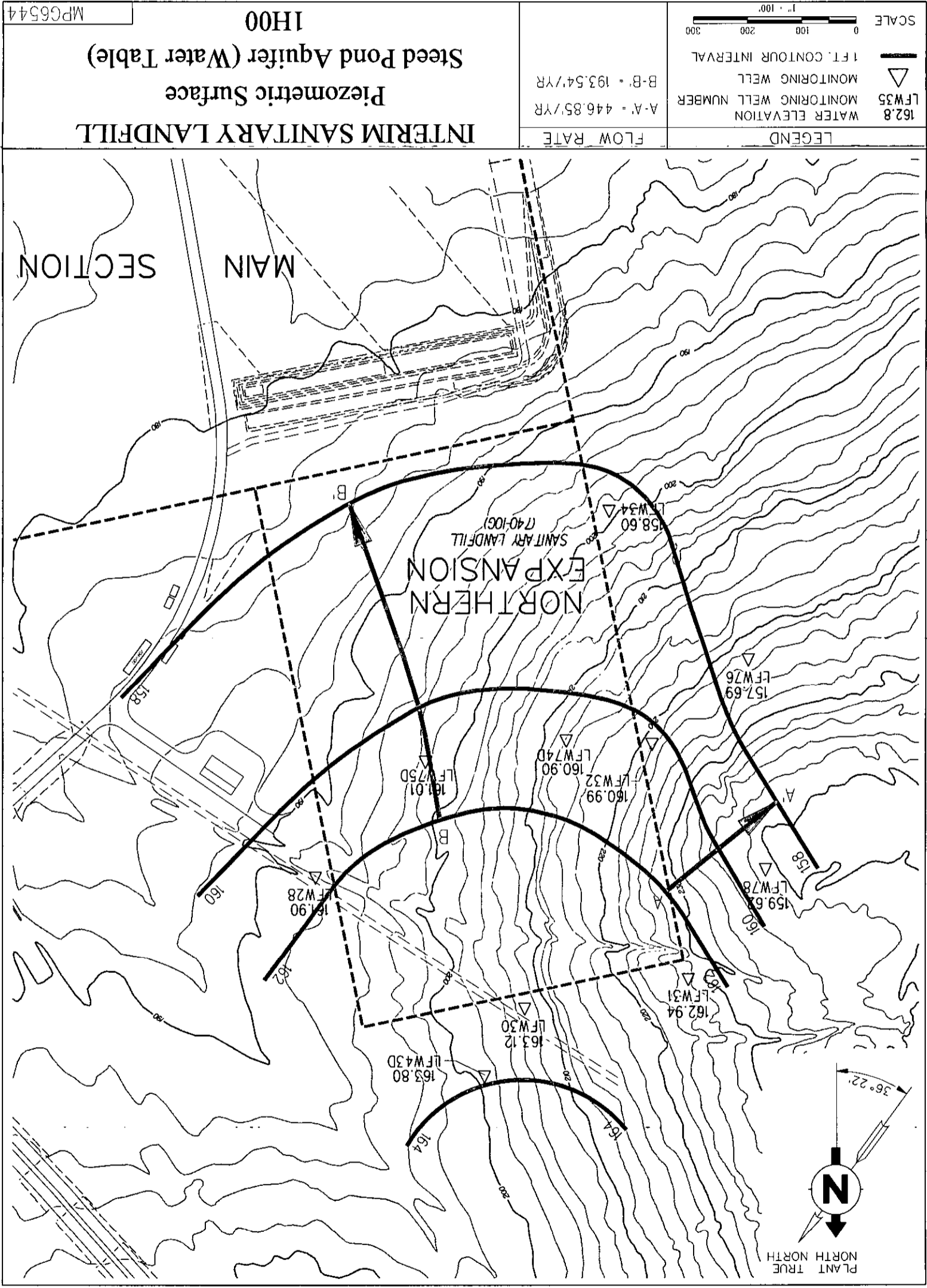


Figure 3. Potentiometric surface map of the Steed Pond Aquifer, 1H00

44434
86826

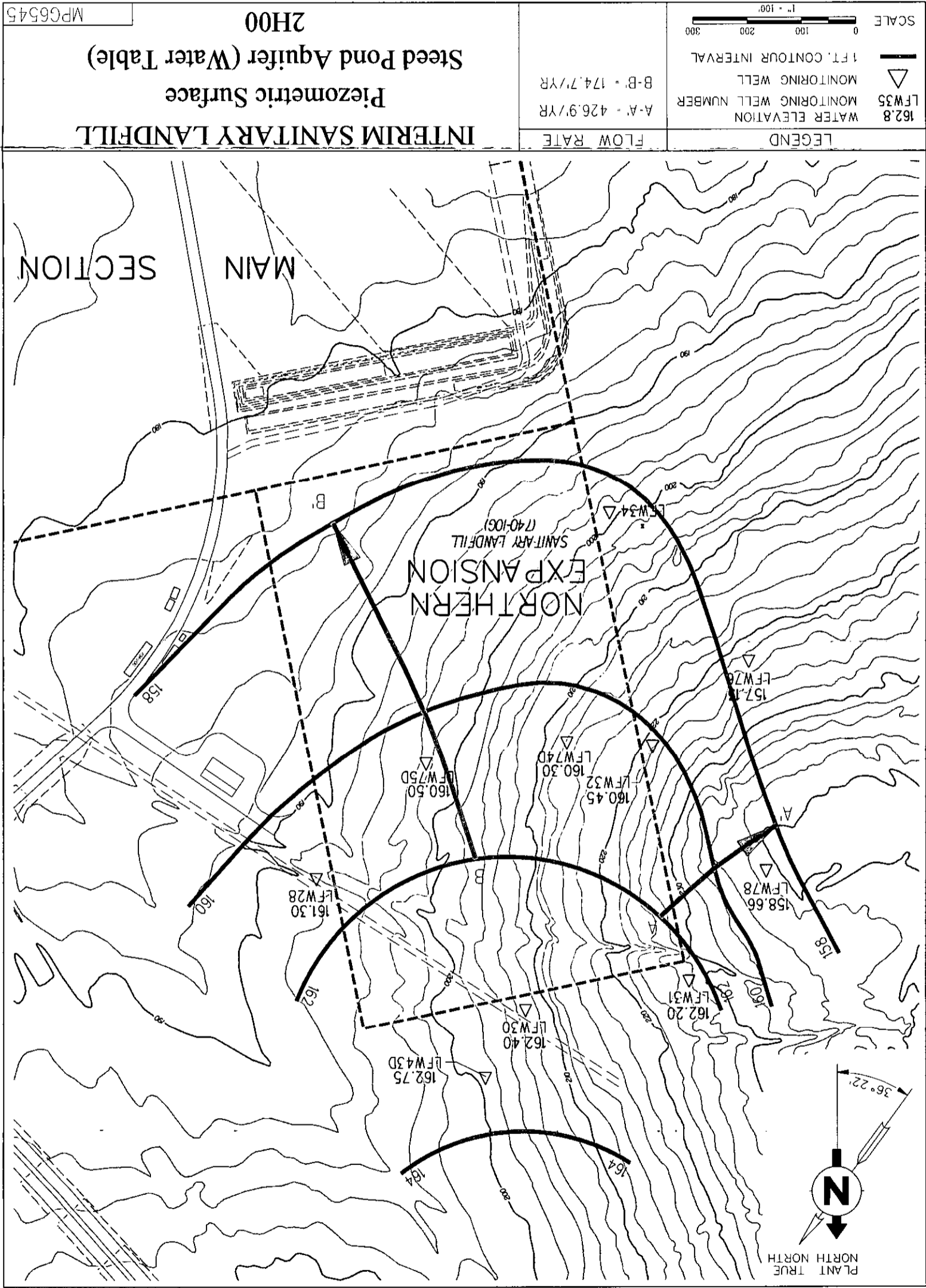


Figure 4. Potentiometric surface map of the Steed Pond Aquifer, 2H00

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N86826-

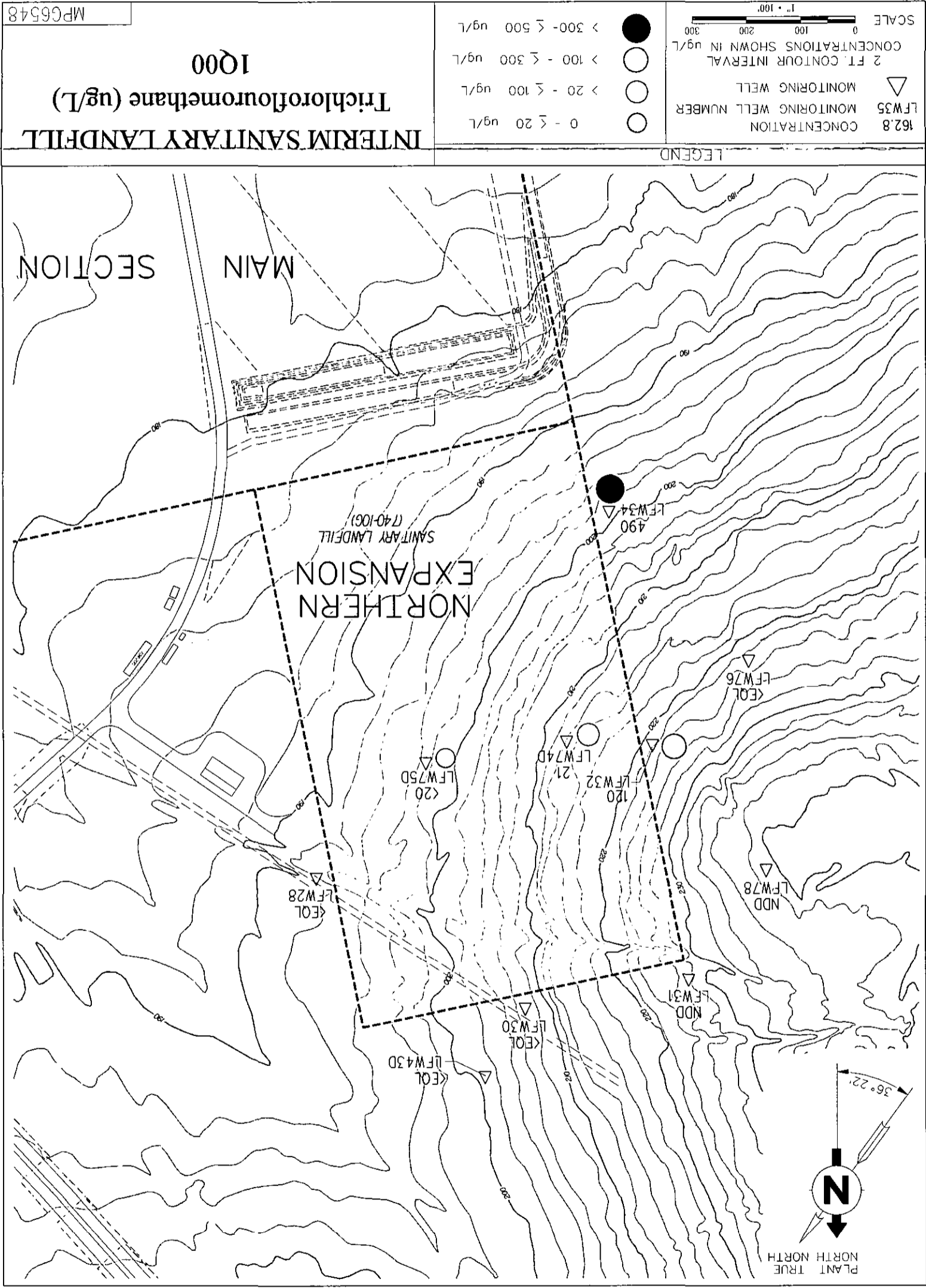


Figure 5. Trichlorofluoromethane (ug/L), 1000

- E44434
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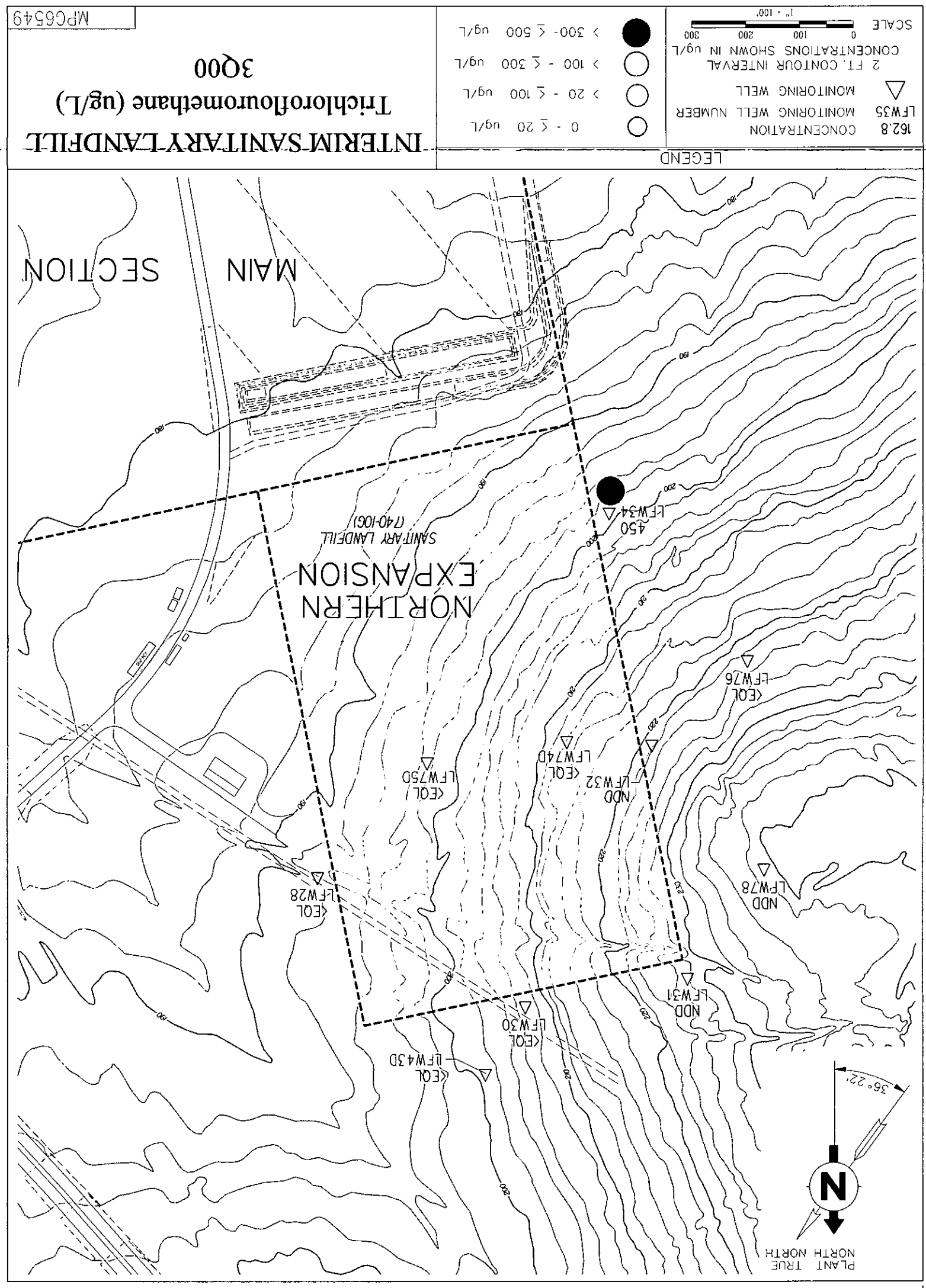


Figure 6. Trichlorofluoromethane (ug/L), 3000

Appendix D

Vadose Methane Monitoring Data

Methane results:

On 10/7/99 4:39 p.m., ADS Laboratory (SRTC/SRS) completed analysis of methane samples submitted the morning of 10/7/99. Analyses were performed by GC-Mass Spec. A 30 m x 0.53 mm Carboxen-1004 porous layer open tubular capillary column was used for the separation.

Table D.1 Methane Monitoring Results

| <u>Lab ID#</u> | <u>Station ID</u> | <u>Methane, volume% (1)</u> |
|----------------|-----------------------|-----------------------------|
| 3-134091 | LGM-1 | 12.1 vol% |
| 3-134092 | LGM-2 | 0.51 vol% |
| 3-134093 | LGM-3 | 3.40 vol% |
| 3-134094 | LGM-4 | 48.5 vol% |
| 3-134094 | LGM-4 (w/5x dilution) | 53.7 vol% * |

(1) Results are reported with three significant figures. The last digit is insignificant, but reported

* Includes dilution factor correction to base value on original sample.

Post calibration verification, after the LGM-4 sample was performed to check instrument for calibration drift. The instrument remained in specification for calibration. The 53.7 vol% value is more reliable since it is within the calibration linear range.

The LEL of methane is ~5%. Hence, any result greater than 5% is > LEL.